

Golder Associates Inc.

305 Fellowship Road, Suite 200
Mt. Laurel, NJ USA 08054
Tel: (609) 273-1110
Fax (609) 273-0778



REVISED DRAFT

**REMEDIAL INVESTIGATION REPORT
BERKS LANDFILL
BERKS COUNTY, PENNSYLVANIA**

VOLUME 3 OF 3

**Appendices N through W
and
Attachments 1 and 2**

Prepared by:

**Golder Associates Inc.
305 Fellowship Road, Suite 200
Mt. Laurel, NJ 08054**

DISTRIBUTION:

4 Copies	-	A. Koller, USEPA
1 Copy	-	R. Klinikowski, PADER
1 Copy	-	R. Bishop, PADER
1 Copy	-	M. Rosenberg, CH2M Hill
5 Copies	-	Berks Landfill Respondents
2 Copies	-	Golder Associates Inc.

March 1995

Project No.: 913-6773

AR302647

TABLE OF CONTENTS**VOLUME 3****LIST OF APPENDICES****Appendix N Aquatic Habitat Assessment****Appendix O Agency Correspondence for NEPA Issues****Appendix P Laboratory Analytical Results
Attachment P-1 Fourth Quarter Residential Well Data****Appendix Q Detected Constituent Summary Tables****Appendix R Procedures to Calculate Summary Statistics of Detected Constituents****Appendix S Monitoring Well C-3D Groundwater Sample Analysis Results****Appendix T Stastical Comparison of Metal Concentrations in Background and
On-Site Groundwater****Appendix U Phase 1B Sediment Sampling Grain Size Distribution Results****Appendix V Results of Phase 1A Air Monitoring****Appendix W Phase 1B EPA Method TO14 Results and Field GC Chromatograms****LIST OF ATTACHMENTS**

Attachment 1	Description of Drum Investigation
Attachment 1-A	Geophysical Investigation Report
Attachment 1-B	Test Pit/Trench Logs
Attachment 1-C	Bill of Loading and Manifests for Material Transportation
Attachment 1-D	Soil Sample Analytical Results

Attachment 2 USEPA Split Sampling Results

APPENDIX N
AQUATIC HABITAT ASSESSMENT

AR302649

APPENDIX N

AQUATIC HABITAT ASSESSMENT

1.0 INTRODUCTION

A benthic macroinvertebrate study was conducted at the Berks Landfill, Spring Township, Berks County, Pennsylvania, as part of the implementation of the Berks Landfill Remedial Investigation/Feasibility Study Work Plan approved by USEPA on June 8, 1992. The purpose of the study was to characterize the composition and structure of the benthic macroinvertebrate communities in streams within and adjacent to the Berks Landfill Site and to assess potential impacts of the Site on aquatic habitats. Characterization of benthic macroinvertebrate communities in streams was accomplished by using qualitative benthos sampling and identification methods. Samples were collected from the tributary to Cacoosing Creek and other unnamed tributaries at the Site. Potential impacts of the Site to aquatic habitats (impairment) were assessed by comparing macroinvertebrate communities upstream and downstream of the Site.

2.0 BENTHIC MACROINVERTEBRATES AS INDICATORS OF ENVIRONMENTAL IMPACTS

Benthic macroinvertebrates (benthos) are aquatic organisms which inhabit the bottom or substratum of streams, estuaries, and other waterbodies. Although some immature benthos (early instars) of many species are extremely small, the majority of benthos are visible to the unaided eye. Benthos serve as the primary food source for many fish and other aquatic animals.

The composition and structure of the benthic communities are generally stable from year to year in undisturbed habitats. Seasonal variations, however, may exist due to life-cycle dynamics. A diverse and balanced distribution of benthic

communities are indicative of most free-flowing aquatic environments that have not been adversely impacted.

Benthos generally have limited ranges and are, therefore, good indicators of localized aquatic conditions. When macroinvertebrate communities are subjected to even minor habitat disturbances (i.e., water quality, siltation, organic load, etc.) they generally respond by adjustments in the community structure. Benthic macroinvertebrate surveys can therefore provide useful information to characterize potential adverse impacts to aquatic habitats.

3.0 METHODS

Rapid Bioassessment Protocol I (RBP I) methodologies, outlined in the Rapid Bioassessment Protocols For Use In Stream And Rivers: Benthic Macroinvertebrates And Fish (EPA/444/4-89-001), 1989 were utilized to evaluate benthic macroinvertebrate communities at the Site. RBP I focuses on the qualitative assessment of benthic communities. It should be noted that Golder conducted additional investigative work beyond the typical RBP I protocol which was not specified in the Work Plan. This additional work was performed to provide a more comprehensive evaluation of the aquatic habitats. Typical RBP III protocols including collection, preservation, and subsequent identification and comparisons of benthic macroinvertebrate specimens, grain size distribution for benthic macroinvertebrate colonization, and more detailed sample station evaluations were conducted. Details of the sample collection and evaluation methods utilized are discussed below. In addition, field measurements of surface water quality and laboratory analyses of surface water and sediment samples were performed. These sampling and analysis procedures are also described below.

3.1 Sample Station Locations

Six sampling stations were established in ecologically similar settings to provide a representative comparison of benthic macroinvertebrate communities. Prior to the selection of the sampling station locations, a review of stream habitats and physical conditions was conducted through stream reconnaissance at the general location of the sampling points identified in the USEPA approved RI/FS Work Plan. Final sample station locations were then selected considering stream substrate, vegetation, water depth, and other stream conditions such as riffle-pool and riffle-run ratios and sun exposure. All the stations were considered to have ecologically similar habitats even though there were certain physical differences.

For future reference and relocation, a marker (placard on a wooden stake) was placed in the general location of each of the sampling stations. Sampling stations were identified as Macroinvertebrate (MI) Station 1 through 6. The general location of the benthic macroinvertebrate sampling stations are shown on Figure N-1.

Two reference or background sampling stations, Station 4 and Station 6, (see Figure N-1) were established. Station 6 was selected upstream of the Site within the unnamed tributary to Cacoosing Creek. Station 4 was selected in the upstream headwaters of the central drainage channel which flows between the western and eastern landfills at the Site. Both stations were selected and sampled in the same manner as all other sampling locations. These reference locations were used for the final comparison of benthos within the Site.

3.2 Sampling Methods

Benthic macroinvertebrate samples were collected using a rectangular aquatic kick net with a mesh size of 800 microns. The net was firmly placed against the stream bottom while the substrata immediately upstream from the net was disturbed,

allowing the benthic macroinvertebrate organisms and debris to be washed into the net by the stream flow. An area of approximately 1-foot by 3-feet was kicked or disturbed by hand for each sample collected. Large rocks within the sample zone, were also examined for the presence of benthos. Eight individual kick samples (collection points), four in the riffle habitat and four in the pool or run habitat, were collected at each of the six sample stations.

The contents of the net were emptied into a sorting tray containing about 1/2 inch of water. The net was then carefully checked for any remaining benthos. Benthic macroinvertebrate organisms were hand picked (with curved forceps; eye droppers; etc.) from the tray and preserved in 2 ounce glass bottles (with polyethylene inserts) using a 10% Formalin solution buffered for tissue fixation. Organic debris and substrata were sorted from all samples prior to preservation. Samples from each collection point at a sample station were placed in separate sample bottles (i.e., 8 bottles for each sample station). It should be noted that RBP I does not require the collection of specimens and that Golder retained the specimens for a more accurate evaluation of the benthos present at the Site. It is also worth noting that, in most cases, greater than 100 organism counts were found at each collection point.

Samples were collected on May 19 through 21, 1992. Representatives from the USEPA and Tetra Tech, Inc. (the USEPA oversight contractor) were present for a portion of the benthic macroinvertebrate sampling. Each collection point was assigned a unique sample identification number including the sample station identification number. All benthic sampling was supplemented with surface water quality field measurements for dissolved oxygen, temperature, pH, and specific conductivity, stream flow measurements, and physical observations and notes. Surface water samples were collected for laboratory analyses of total suspended solids, alkalinity and hardness. In addition, sediment samples were collected for laboratory analyses (i.e., grain size distribution, total organic carbon, percent

moisture, percent solids). Field measurements, sample collection and laboratory analyses were conducted in accordance with the USEPA approved RI/FS Work Plan.

3.3 Evaluation of Benthic Macroinvertebrate Samples

Benthic macroinvertebrate organism samples were examined with an illuminated stereomicroscope. Identifications were conducted to an Order and/or Family taxonomic level (refer to Table N-1). Sample specimens were then compared with the background specimens collected at the Site and with known pollutant sensitive benthic macroinvertebrate groups (established by USEPA) to evaluate stream environmental conditions.

4.0 MACROINVERTEBRATE SURVEY RESULTS

Complexity in benthic macroinvertebrate community structure is mainly determined by the numbers of individuals present, the number of taxa (species) present and the type and distribution of the taxa found within the community. Density, taxa richness, and species diversity are descriptive parameters used to evaluate these various community characteristics. The Biosurvey Field Data Sheets (provided in Attachment N-1) present the results of the benthic macroinvertebrate surveys for the Berks Landfill Site. These data are intended to reflect the relative complexity of the community structures and provide a comparison of each of the communities observed.

The results indicate that, within the benthic macroinvertebrate communities assessed both upstream and downstream of the Site, there is an overall benthic macroinvertebrate abundance and rich taxa diversity. There is also a dominance of the following pollutant sensitive or intolerant organisms, as identified by USEPA: Ephemeroptera (mayflies), Plecoptera (stoneflies) and Trichoptera

(caddisflies) collectively referred to as "EPT". In addition, there is a relatively low distribution of pollutant tolerant groups such as Diptera and Annelida (e.g., Chironomidae and Oligochaetes).

Minnows (Notropis sp.) were observed at five (5) of the sampling stations. Station 4 was the only station where minnows were not observed. It should be noted, however, that sampling Station 4 is an upstream location and is within the first order stream headwaters of the central drainage channel. Fish would typically not be expected within these areas.

The benthic macroinvertebrate survey results indicate that a healthy aquatic habitat exists within the streams sampled and that there is a dominance of EPT species which indicate that suitable conditions exist within the stream systems at the Site to support normal benthic communities. Distribution of species appeared to be consistent throughout the stream system and overall species diversity was also observed which reflects the similarity of conditions at each of the individual sampling stations.

Water quality parameters measured in the field included pH, temperature, specific conductance and dissolved oxygen. The results are provided in Table N-2. Chemical analyses of surface water samples for alkalinity, hardness, and Total Suspended Solids are provided in Table N-3. The range of field parameter values measured and analytical results for the conventional parameters sampled during this study were as follows:

pH	6.46	-	7.90	standard units;
specific conductance	220	-	330	umhos/cm;
dissolved oxygen	7.9	-	9.7	mg/l;
temperature	10.6°C	-	17.7°C;	
alkalinity	72	-	126	mg/l as CaCO ₃ ;
hardness	137	-	188	mg/l as CaCO ₃ ; and
total suspended solids	<xx	-	23	mg/l.

The Pennsylvania Water Quality Standards, outlined in Pennsylvania Code Title 25, Chapter 93 define the designated use of Cacoosing Creek as a Warm Water Fishery (WWF). The WWF designated use criteria established for the measured parameters are as follows:

pH	6.0 - 9.0 standard units;
total dissolved solids (specific conductance)	1500 mg/l maximum (specific conductance x 0.65 = TDS);
dissolved oxygen	minimum daily average 5.0 mg/l, any one minimum 4.0 mg/l;
temperature	maximum temperature 22.2°C for May 16 through May 31;
alkalinity	minimum of 20 mg/l as CaCO ₃ ;
hardness	maximum monthly average 150 mg/l as CaCO ₃ ; and
total suspended solids	no criteria established.

All of the values, except hardness, were within the PADER criteria for the designated use of the surface water body. It should be noted that, while three of the six sample locations minimally exceeded the hardness criteria, one of the three samples was from a reference location (Station 6) which may indicate that the hardness levels measured are indicative of natural stream conditions.

Table N-4 presents the inorganic chemical analyses results for Total Organic Carbon (TOC) and percent solids for stream sediments. These results indicate that TOC concentrations are considerably low (i.e., near the detection limit of the analytical method). These results also indicate suitable conditions for benthic macroinvertebrate communities.

The results of the grain size distribution analyses (summary provided in Table N-5 and in Attachment N-2 for laboratory data) show that stream habitat conditions, with regard to substrate characteristics, are relatively similar and that colonization within benthic communities could therefore be expected to be similar at each of

the sampling station locations. It should be noted the grain size distribution conducted during the benthos survey was to evaluate benthic macroinvertebrate colonization habitat similarities.

Stream flow measurements were also performed at each macroinvertebrate station. Stream flow were measured by the standard USGS Streamflow Measurement Technique using a pygmy current meter suspended on a top setting wading rod. Results of the measurements are presented in Table N-6 and cross sections of the streams are presented on Figure N-2.

5.0 SUMMARY OF RESULTS

The results of the benthic macroinvertebrate survey, field measured surface water quality parameters and laboratory analyses of surface water samples for conventional parameters indicate that the Berks Landfill Site has not adversely impacted the benthic community structure and, has not exceeded the criteria to protect the designated use of the streams. This conclusion is evident by: (1) the diverse benthic community structure with a high occurrence of taxa present; (2) high population densities at each benthic sampling station; (3) a dominance of EPT (pollutant intolerant) species which are abundant and generally widespread throughout the stream systems at the Site; (4) a low occurrence of pollutant tolerant species such as those in the Order Diptera; and, (5) surface water parameters measured in the field and the laboratory generally meet PADER surface water quality criteria for the designated use of the stream. In summary, the data collected indicates that healthy aquatic habitats exist at the Berks Landfill Site.

REFERENCES

- Adler, P.H. and K.C. Kim. February 1986. The Blackflies of Pennsylvania (Diptera), Bionomics, Taxonomy, and Distribution. The Pennsylvania State University, College of Agriculture, Agricultural Experiment Station, Bulletin 856., University Park, Pennsylvania. 88 pp.
- Klemm, Donald J. et al. November 1990. Macroinvertebrate Field and Laboratory Methods for Evaluating the Biological Integrity of Surface Waters. United States Environmental Protection Agency. Cincinnati, Ohio EPA/600/4-90/030.
- Merritt, R.W., and K.W. Cummins. 1984. An Introduction to the Aquatic Insects of North America, Second Edition. Kendall/Hunt Publishing Co., Dubuque, Iowa. 721 pp.
- PADER. 1988. Guidelines for Benthic Macroinvertebrates Stream Surveys for Landfills, June 1988. PADER Bureau of Waste Management, Harrisburg, PA.
- Plafkin, J.L., et al. May 1989. Rapid Bioassessment Protocols For Use in Streams and Rivers, Benthic Macroinvertebrates and Fish. United States Environmental Protection Agency. Washington, D.C. EPA/444/4-89-001.
- Surdick, R.F., and K.C. Kim. May 1976. Stoneflies (Plecoptera) of Pennsylvania. The Pennsylvania State University, College of Agriculture, 125 Agricultural Experiment Station, Bulletin 808., University Park, Pennsylvania. 73 pp.
- Thorp, J.H. and A.P. Covich. 1991. Ecology and Classification of North American Freshwater Invertebrates. Academic Press, Inc. San Diego, California. 910 pp.

TABLE N-1

LIST OF MACROINVERTEBRATES OBSERVED AT THE BERKS LANDFILL

ORDER

Suborder
Family
Genera

EPHEMEROPTERA

Suborder Pannota
Ephemerellidae
Suborder Schistonota
Baetidae
Heptageniidae
Leptophlebiidae

PLECOPTERA

Suborder Systelognatha
Chloroperlidae
Perlidae
Perlodidae
Suborder Euholognatha
Leuctridae
Nemouridae

TRICHOPTERA

Suborder Annulipalpia
Hydropsychidae
Psychomyiidae
Suborder Spicipalpia
Hydroptilidae
Suborder Integripalpia
Limnephilidae
Uenoidae

DIPTERA

Chironomidae
Empididae
Simuliidae
Suborder Nematocera
Tipulidae
Tipula sp.

COLEOPTERA

Elmidae
Microcyloepus sp.
Heterimnius sp.
Psephenidae

HEMIPTERA

Gerridae
Gerris sp.
Valiidae
Microvelia sp.

DECAPODA

Cambaridae
Cambrus sp.

ODONATA

Suborder Anisoptera
Gomphidae

TABLE N-2
 RESULTS OF SURFACE WATER SAMPLES
 BERKS LANDFILL REMEDIAL INVESTIGATION - MAY 1992
 BERKS COUNTY, PENNSYLVANIA

Parameter	Units	Primary Sample Results FIELD PARAMETERS					
		Sample Location MI STATION 1 Date Sampled: 05/19/92	Sample Location MI STATION 2 Date Sampled: 05/19/92	Sample Location MI STATION 3 Date Sampled: 05/20/92	Sample Location MI STATION 4 Date Sampled: 05/20/92	Sample Location MI STATION 5 Date Sampled: 05/20/92	Sample Location MI STATION 6 Date Sampled: 05/21/92
		Result	Result	Result	Result	Result	Result
pH	std. units	7.64	7.90	6.46	7.20	7.51	7.02
Specific Conductance	umhos/cm	330	320	285	220	270	240
Dissolved Oxygen	mg/l	9.2	7.9	9.6	9.1	8.5	9.7
Temperature	Deg. C	13.2	14.2	10.6	13.5	17.7	11.1
Temperature using traceable thermometer	Deg. C	13.5	15.2	10.8	13.5	17.5	11.2

Notes:

1. MI Indicates Macroinvertebrate
2. MI STATION 1 is SW-08 monitoring location as shown on Figure 2 of the FSP.
3. MI STATION 2 is SW-09 monitoring location as shown on Figure 2 of the FSP.
4. MI STATION 3 is SW-10 monitoring location as shown on Figure 2 of the FSP.
5. MI STATION 4 is SW-04 monitoring location as shown on Figure 2 of the FSP.
6. MI STATION 5 is SW-11 monitoring location as shown on Figure 2 of the FSP.
7. MI STATION 6 is SW-01 monitoring location as shown on Figure 2 of the FSP.

TABLE N-3
 RESULTS OF SURFACE WATER SAMPLES
 BERKS LANDFILL REMEDIAL INVESTIGATION - MAY 1992
 BERKS COUNTY, PENNSYLVANIA

Parameter	Primary Sample Results INORGANICS																	
	Sample Location SW-01/1A Date Sampled: 05/21/92			Sample Location SW-04/1A Date Sampled: 05/20/92			Sample Location SW-08/1A Date Sampled: 05/19/92			Sample Location SW-09/1A Date Sampled: 05/19/92			Sample Location SW-10/1A Date Sampled: 05/20/92			Sample Location SW-11/1A Date Sampled: 05/20/92		
	MRL	Result	Qual	MRL	Result	Qual	MRL	Result	Qual	MRL	Result	Qual	MRL	Result	Qual	MRL	Result	Qual
Alkalinity	2	92	A	2	72	A	2	128	A	2	120	A	2	92	A	2	120	A
Hardness	2	151	A	2	137	A	2	188	A	2	163	A	2	144	A	2	139	A
Total Suspended Solids	5	13	A	5	ND	U	5	ND	U	5	23	A	5	20	A	5	23	A

- Notes:**
- All results are reported in units of mg/l.
 - MRL is the laboratory method reporting limit.
 - ND indicates that this parameter was not detected in this sample.
 - Qualifiers (Qual) are indicated as follows:
 A - Acceptable (Quantitative) Data
 J - Estimated (Semi-quantitative) Data
 K - Estimated Biased High Data
 L - Estimated, Biased Low Data
 R - Unusable Data
 U - Not Detected (Quantitative) Data
 UJ - Not Detected (Semi-quantitative) Data
 UL - Not Detected, Biased Low Data
 - Sample location SW-01 is Macroinvertebrate Station 6.
 Sample location SW-04 is Macroinvertebrate Station 4.
 Sample location SW-08 is Macroinvertebrate Station 1.
 Sample location SW-09 is Macroinvertebrate Station 2.
 Sample location SW-10 is Macroinvertebrate Station 3.
 Sample location SW-11 is Macroinvertebrate Station 5.

AR302661

TABLEN-4

RESULTS OF SEDIMENT SAMPLES

BERKS LANDFILL REMEDIAL INVESTIGATION - MAY 1992
BERKS COUNTY, PENNSYLVANIA

Parameter	Sample Location SED01/1A Date Sampled: 05/21/92		Sample Location SED03/1A Date Sampled: 05/20/92		Sample Location SED07/1A Date Sampled: 05/20/92		Sample Location SED08/1A Date Sampled: 05/19/92		
	MRL	Result	Qual	MRL	Result	Qual	MRL	Result	Qual
Total Organic Carbon (TOC)	0.1	0.4	A	0.1	0.5	A	0.1	0.2	A
Percent Solid	NA	76	NA	NA	76	NA	NA	73	NA

NOTES:

- All sediment results are reported as a percent and the Rinsale Blank results are reported in mg/L. TOC results are reported on a dry weight basis.
- MRL is the laboratory method reporting limit.
- ND indicates that this parameter was not detected in this sample.
- NA indicates Not Applicable.
- Qualifiers (Qual) are indicated as follows:
 A - Acceptable (Quantitative) Data
 J - Estimated (Semi-quantitative) Data
 UJ - Not Detected (Semi-quantitative) Data
 UL - Not Detected, Biased Low Data
 R - Unusable Data
 L - Estimated, Biased Low Data
 K - Estimated Biased High Data
 U - Not Detected (Quantitative) Data
- Sample Location SED01/1A is Macroinvertebrate Station 6.
 Sample Location SED03/1A is Macroinvertebrate Station 4.
 Sample Location SED07/1A is Macroinvertebrate Station 5.
 Sample Location SED08/1A is Macroinvertebrate Station 1.
 Sample Location SED09/1A is Macroinvertebrate Station 3.
 Sample Location SED10/1A is Macroinvertebrate Station 2.
 RB001/1A sample is the Rinsale Blank Sample.

AR302662

TABLE N-4 (Cont'd)

RESULTS OF SEDIMENT SAMPLES

BERKS LANDFILL REMEDIAL INVESTIGATION - MAY 1992
BERKS COUNTY, PENNSYLVANIA

Parameter	Primary Sample Results INORGANICS											
	Sample Location SED09/1A			Sample Location SED10/1A			Sample Location RB001/1A					
	MRL	Result	Qual	MRL	Result	Qual	MRL	Result	Qual	MRL	Result	Qual
Total Organic Carbon (TOC)	0.1	0.3	A	0.1	0.4	A	1	1	ND	1	ND	U
Percent Solid	NA	83	NA	NA	74	NA	NA	NA	NA	NA	NA	NA

Notes:

- All sediment results are reported as a percent and the Flinsate Blank results are reported in mg/l. TOC results are reported on a dry weight basis.
- MRL is the laboratory method reporting limit.
- ND indicates that this parameter was not detected in this sample.
- NA indicates Not Applicable.
- Qualifiers (Qual) are indicated as follows:
 A - Acceptable (Quantitative) Data
 J - Estimated (Semi-quantitative) Data
 UJ - Not Detected (Semi-quantitative) Data
 UL - Not Detected, Biased Low Data
 R - Unusable Data
 L - Estimated, Biased Low Data
 K - Estimated Biased High Data
 U - Not Detected (Quantitative) Data
- Sample Location SED01/1A is Macroinvertebrate Station 6.
 Sample Location SED03/1A is Macroinvertebrate Station 4.
 Sample Location SED07/1A is Macroinvertebrate Station 5.
 Sample Location SED08/1A is Macroinvertebrate Station 1.
 Sample Location SED09/1A is Macroinvertebrate Station 3.
 Sample Location SED10/1A is Macroinvertebrate Station 2.
 RB001/1A sample is the Flinsate Blank Sample.

AR302663

TABLE N-5
SUMMARY OF RESULTS OF GRAIN SIZE DISTRIBUTION
BERKS LANDFILL REMEDIAL INVESTIGATION - MAY 1992
BERKS COUNTY, PENNSYLVANIA

Macro-Invertebrate Station	Sample Number	Delivered Moisture Content	Distribution		
			% Finer #4 Sieve	% Finer #200 Sieve	% Finer 2 Micron
1	GSD-1	11.1%	31.4%	1.6%	0.3%
2	GSD-2	21.5%	45.5%	4.9%	0.7%
3	GSD-3	18.4%	23.6%	2.0%	0.2%
4	GSD-4	21.7%	42.3%	1.7%	0.0%
5	GSD-5	21.4%	51.8%	2.0%	0.5%
6	GSD-6	32.0%	65.3%	6.2%	1.6%

AR302664

TABLE N-6
STREAM FLOW MEASUREMENTS
BERKS LANDFILL REMEDIAL INVESTIGATION -- MAY 1992
BERKS COUNTY, PENNSYLVANIA

STATION ID	LOCATION	DATE	METHOD	FLOW GPM
MI-1	Macroinvertebrate Station 1	5/18/92	Pygmy Meter	386.0
MI-2	Macroinvertebrate Station 2	5/19/92	Pygmy Meter	52.5
MI-3	Macroinvertebrate Station 3	5/20/92	Pygmy Meter	104.1
MI-4	Macroinvertebrate Station 4	5/20/92	Pygmy Meter	34.6
MI-5	Macroinvertebrate Station 5	5/18/92	Pygmy Meter	187.5
MI-6	Macroinvertebrate Station 6	5/21/92	Pygmy Meter	84.4

AR302665

LABORATORY DATA QUALIFIER LEGEND

ORGANICS

- U - Indicates compound was analyzed for but not detected.
- J - Indicates an estimated value.
- B - This flag is used when the analyte is found in the associated blank as well as in the sample. It indicates possible/probable blank contamination and warns the data user to take appropriate action.

METALS AND CYANIDES

- E - The reported value is estimated because of the presence of interference.
- N - Spiked sample recovery not within control limits.
- W - Post-digestion spike for Furnace AA analysis is out of control limits (85-115%), while sample absorbance is less than 50% of spike absorbance.
- * - Duplicate analysis not within control limits.

AR302666



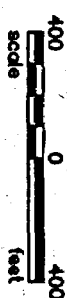
REV	DATE	BY	DESCRIPTION
01	03/25/94	AS SHOWN	
02	03/25/94	AS SHOWN	
03	03/25/94	AS SHOWN	
04	03/25/94	AS SHOWN	
05	03/25/94	AS SHOWN	
06	03/25/94	AS SHOWN	
07	03/25/94	AS SHOWN	
08	03/25/94	AS SHOWN	
09	03/25/94	AS SHOWN	
10	03/25/94	AS SHOWN	
11	03/25/94	AS SHOWN	

Golder Associates

BERKS LANDFILL

N-1

**BENTHIC MACROINVERTEBRATE
SAMPLE LOCATIONS**



MAR 28 1994

AR302667

LEGEND

- BENTHIC MACROINVERTEBRATE SAMPLE LOCATIONS
- APPROXIMATE SITE BOUNDARY

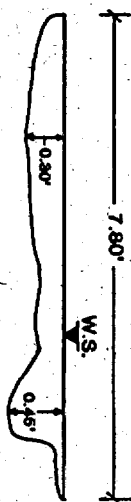
NOTES

- 1) SOME OF THE MACROINVERTEBRATE SAMPLING LOCATIONS ARE FIELD LOCATED.

REFERENCE

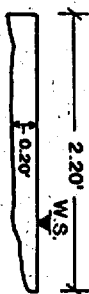
- 1) TOPOGRAPHIC MAPPING PROVIDED BY GEO COOPERATION, NEWQUANDLAND, NJ AERIAL PHOTOGRAPHY DATE DECEMBER 10, 1991.
- 2) HORIZONTAL COORDINATE SYSTEM IS PENNSYLVANIA STATE PLANE NORTH AMERICAN DATUM 1983 AND VERTICAL DATA ARE TIED TO MGD 1928.
- 3) SITE BOUNDARY TAKEN FROM DRAWINGS PREPARED BY L. ROBERT KIMBALL & ASSOCIATES, CONSULTING ENGINEERS & ARCHITECTS, EBERSBURG, PA. MAP NOS. 4375 02 AND 4375 04 TITLED "PROPERTY MAP OF BERKS COUNTY PENNSYLVANIA", DATED NOVEMBER 05, 1983. PROJECTED FROM PA STATE PLANE NA027 TO NA083 BY GOLDER ASSOCIATES USING ARCAD GEOGRAPHIC INFORMATION SYSTEM.

**MI-1
MACROINVERTEBRATE
STATION 1**



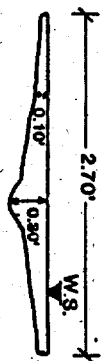
VIEW: DOWNSTREAM
BOTTOM: COBBLES, GRAVELS AND SAND
AREA: 1.97 ft²
VELOCITY: 0.43 ft/sec
DISCHARGE: 386 gpm

**MI-2
MACROINVERTEBRATE
STATION 2**



VIEW: DOWNSTREAM
BOTTOM: SAND AND GRAVEL
AREA: 0.38 ft²
VELOCITY: 0.31 ft/sec
DISCHARGE: 62.5 gpm

**MI-3
MACROINVERTEBRATE
STATION 3**



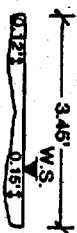
VIEW: DOWNSTREAM
BOTTOM: COBBLES, SAND AND GRAVEL
AREA: 0.36 ft²
VELOCITY: 0.66 ft/sec
DISCHARGE: 104.1 gpm

**MI-4
MACROINVERTEBRATE
STATION 4**



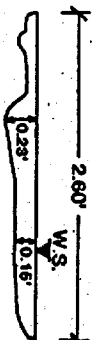
VIEW: DOWNSTREAM
BOTTOM: SAND AND GRAVEL
AREA: 0.16 ft²
VELOCITY: 0.51 ft/sec
DISCHARGE: 34.6 gpm

**MI-5
MACROINVERTEBRATE
STATION 5**



VIEW: DOWNSTREAM
BOTTOM: GRAVEL
AREA: 0.36 ft²
VELOCITY: 1.17 ft/sec
DISCHARGE: 187.5 gpm

**MI-6
MACROINVERTEBRATE
STATION 6**



VIEW: DOWNSTREAM
BOTTOM: SAND AND GRAVEL
AREA: 0.37 ft²
VELOCITY: 0.51 ft/sec
DISCHARGE: 84.4 gpm

MAR 28 1994

AR302668

JOB NO.	913-8773	SCALE	N.T.S.
DR. BY	MRM	DATE	12/19/93
CHE. BY	KCC	FILE NO.	PA21-394
REV. BY	KSW	NO. SHEETS	11

Goldier Associates

STREAM CROSS SECTIONS
MAY 1992 SAMPLING EVENT

BERKS LANDFILL

DATE: N-2

Attachment N-1
Biosurvey Data Forms

AR302669

Benthic Macroinvertebrate Organism Sampling

MI-1
SAMPLE STATION ID.

MAY 19, 1992
DATE

Rapid Bioassessment Protocol I

Biosurvey Field Data Sheet

RELATIVE ABUNDANCE OF AQUATIC BIOTA

Periphyton	0	1	2	3	4		Slimes	0	1	2	3	4
Filamentous Algae	0	①	2	3	4		Macroinvertebrates	0	1	2	3	④
Macrophytes	0	1	2	3	4		Fish	0	1	2	③	4

0 = Absent/Not Observed 1 = Rare 2 = Common 3 = Abundant 4 = Dominant

MACROBENTHOS QUALITATIVE SAMPLE LIST (Indicate Relative Abundance R = Rare, C = Common, A = Abundant, D = Dominant)

Porifera	Anisoptera	Chironomidae	R
Hydrozoa	Zygoptera	Plecoptera	D
Pisthnelminthes	Hemiptera	Ephemeroptera	D
Turbellaria	Coleoptera	Trichoptera	R
Hirudinea	Lepidoptera	Other	
Oligochaeta	Sialidae		
Isopoda	Corydalidae		
Amphipoda	Tipulidae		R
Decapoda	Empididae		
Gastropoda	Simuliidae		R
Bivalvia	Taberidae		
	Culicidae		

Rare < 3 Common 3-8 Abundant > 10 Dominant > 50 (Estimate)

Observations

Plant species observed:

- | | |
|---|---|
| <ul style="list-style-type: none"> mayapple Sycamore honeysuckle (stim) grasses Curly dock garlic mustard Creeping buttercup christmas fern | <ul style="list-style-type: none"> multiflora rose cleaverweed wild geranium jack in the pulpit |
|---|---|

GOLDER PROJECT NO. 913-6773

Reference: Figure taken from USEPA Guidance
Manual No. EPA/444/4-89-001, 1989
Figure 6.1-1 pg. 6-2

Benthic Macroinvertebrate Organism Sampling

MI-1
SAMPLE STATION ID.

MAY 19 1992
DATE

IMPAIRMENT ASSESSMENT SHEET

1. Detection of impairment: Impairment detected (Complete items 2-6) **No impairment detected (Stop here)**

2. Biological impairment indicator:

Benthic macroinvertebrates	Other aquatic communities
<input type="checkbox"/> absence of EPT taxa	<input type="checkbox"/> Periphyton
<input type="checkbox"/> dominance of tolerant groups	<input type="checkbox"/> filamentous
<input type="checkbox"/> low benthic abundance	<input type="checkbox"/> other
<input type="checkbox"/> low taxa richness	<input type="checkbox"/> Macrophytes
<input type="checkbox"/> other	<input type="checkbox"/> Slimes
	<input type="checkbox"/> Fish

3. Brief description of problem: _____

Year and date of previous surveys: _____

Survey data available in: _____

4. Cause: (indicate major cause) organic enrichment toxicants flow
habitat limitations other _____

5. Estimated areal extent of problem (m^2) and length of stream reach affected (m), where applicable: _____

6. Suspected source(s) of problem:

point source discharge (name, type of facility, location)
 construction site runoff
 combined sewer outfall
 silviculture runoff
 animal feedlot
 agricultural runoff
 urban runoff
 ground water
 other
 unknown

Briefly explain:

GOLDER PROJECT NO. 913-6773

Reference: Figure taken from USEPA Guidance
Manual No. EPA/444/4-89-001, 1989
Figure 6.1-2 pg. 6-3

Benthic Macroinvertebrate Organism Sampling

MI-2
SAMPLE STATION ID.

MAY 19, 1992
DATE

Rapid Bioassessment Protocol I

Biosurvey Field Data Sheet

RELATIVE ABUNDANCE OF AQUATIC BIOTA

Periphyton	0	1	2	3	4		0	1	2	3	4
Filamentous Algae	0	1	2	3	4	Slimes	0	1	2	3	4
Macrophytes	0	1	2	3	4	Macroinvertebrates	0	1	2	3	4
						Fish	0	1	2	3	4

0 = Absent/Not Observed 1 = Rare 2 = Common 3 = Abundant 4 = Dominant

MACROBENTHOS QUALITATIVE SAMPLE LIST (Indicate Relative Abundance R = Rare, C = Common, A = Abundant, D = Dominant)

Pterinea	Anisoptera	R	Chironomidae	R
Hydrozoa	Zygoptera		Plecoptera	D
Platyhelminthes	Hemiptera	R	Ephemeroptera	D
Turbellaria	Coleoptera	R	Trichoptera	C
Hirudinea	Lepidoptera		Other	
Oligochaeta	Stelidae		Newt: <i>Desmictylus</i> sp.	
Isopoda	Corydalidae			
Amphipoda	Tipulidae	R		
Decapoda	Empididae			
Gastropoda	Simuliidae			
Bivalvia	Tabanidae			
	Culicidae			

Rare < 3 Common 3-8 Abundant > 10 Dominant > 50 (Estimate)

Observations

Plant species observed:

- multiflora rose
- clearweed
- mayapple
- oak
- curly dock
- honey suckle
- shagbark hickory
- sycamore

- christmas fern
- tulip tree
- spice bush
- poison ivy
- wild onion
- wild geranium

- creeping buttercup
- garlic mustard
- jack in the pulpit

GOLDER PROJECT NO. 913-6773

Reference: Figure taken from USEPA Guidance
Manual No. EPA/444/4-89-001, 1989
Figure 6.1-1 pg. 6-2

Benthic Macroinvertebrate Organism Sampling

MI-2
SAMPLE STATION ID.

MAY 19 1993
DATE

IMPAIRMENT ASSESSMENT SHEET

1. Detection of impairment: Impairment detected (Complete items 2-6) **No impairment detected (stop here)**
2. Biological impairment indicator:
- | Benthic macroinvertebrates | Other aquatic communities |
|---|--------------------------------------|
| <input type="checkbox"/> absence of EPT taxa | <input type="checkbox"/> Periphyton |
| <input type="checkbox"/> dominance of tolerant groups | <input type="checkbox"/> filamentous |
| <input type="checkbox"/> low benthic abundance | <input type="checkbox"/> other |
| <input type="checkbox"/> low taxa richness | <input type="checkbox"/> Macrophytes |
| <input type="checkbox"/> other | <input type="checkbox"/> Slimes |
| | <input type="checkbox"/> Fish |
3. Brief description of problem: _____
Year and date of previous surveys: _____
Survey data available in: _____
4. Cause: (indicate major cause) organic enrichment toxicants flow
habitat limitations other _____
5. Estimated areal extent of problem (m²) and length of stream reach affected (m), where applicable: _____
6. Suspected source(s) of problem:
- point source discharge (name, type of facility, location)
 - construction site runoff
 - combined sewer outfall
 - silviculture runoff
 - animal feedlot
 - agricultural runoff
 - urban runoff
 - ground water
 - other
 - unknown

Briefly explain:

GOLDER PROJECT NO. 913-6773

Reference: Figure taken from USEPA Guidance
Manual No. EPA/444/4-89-001, 1989
Figure 6.1-2 pg. 6-3

Benthic Macroinvertebrate Organism Sampling

MI-3
SAMPLE STATION ID.

MAY 20, 1992
DATE

Rapid Bioassessment Protocol I

Biosurvey Field Data Sheet

RELATIVE ABUNDANCE OF AQUATIC BIOTA

Periphyton	0	1	2	3	4		Slimes	0	1	2	3	4
Filamentous Algae	0	①	2	3	4		Macroinvertebrates	0	1	2	③	4
Macrophytes	0	1	2	3	4		Fish	0	1	②	3	4

0 = Absent/Not Observed 1 = Rare 2 = Common 3 = Abundant 4 = Dominant

MACROBENTHOS QUALITATIVE SAMPLE LIST (indicate Relative Abundance R = Rare, C = Common, A = Abundant, D = Dominant)

Porifera	Anisoptera	R	Chironomidae	R
Hydrozoa	Zygoptera		Plecoptera	D
Platyhelminthes	Hemiptera		Ephemeroptera	D
Turbellaria	Coleoptera	R	Trichoptera	R
Nirudinea	Lepidoptera		Other	
Oligochaeta	Sialidae		note: Diemictylus sp.	
Isopoda	Corydalidae			
Amphipoda	Tipulidae	R		
Decapoda	Empididae			
Gastropoda	Simuliidae			
Bivalvia	Tabanidae			
	Culicidae			

Rare < 3 Common 3-8 Abundant > 10 Dominant > 50 (Estimate)

Observations

Plant species observed:
 mayapple
 jack in the pulpit
 honeysuckle
 multiflora rose
 ash
 cleaverweed
 vitis sp.

Poison ivy
 curly dock
 moss
 christmas fern
 ground ivy

Common reed
 garlic mustard

GOLDER PROJECT NO. 913-6773

Reference: Figure taken from USEPA Guidance
 Manual No. EPA/444/4-89-001, 1989
 Figure 6.1-1 pg. 6-2

Benthic Macroinvertebrate Organism Sampling

MI-3
SAMPLE STATION ID.

MAY 20, 1992
DATE

IMPAIRMENT ASSESSMENT SHEET

1. Detection of impairment: Impairment detected (Complete items 2-6) No impairment detected (Stop here)
2. Biological impairment indicator:
- | | |
|---|--------------------------------------|
| Benthic macroinvertebrates | Other aquatic communities |
| <input type="checkbox"/> absence of EPT taxa | <input type="checkbox"/> Periphyton |
| <input type="checkbox"/> dominance of tolerant groups | <input type="checkbox"/> filamentous |
| <input type="checkbox"/> low benthic abundance | <input type="checkbox"/> other |
| <input type="checkbox"/> low taxa richness | <input type="checkbox"/> Macrophytes |
| <input type="checkbox"/> other | <input type="checkbox"/> Slimes |
| | <input type="checkbox"/> Fish |
3. Brief description of problem: _____
Year and date of previous surveys: _____
Survey data available in: _____
4. Cause: (indicate major cause) organic enrichment toxicants flow habitat limitations other _____
5. Estimated areal extent of problem (m²) and length of stream reach affected (m), where applicable: _____
6. Suspected source(s) of problem:
- point source discharge (name, type of facility, location)
 - construction site runoff
 - combined sewer outfall
 - silviculture runoff
 - animal feedlot
 - agricultural runoff
 - urban runoff
 - ground water
 - other
 - unknown

Briefly explain:

GOLDER PROJECT NO. _____

Reference: Figure taken from USEPA Guidance
Manual No. EPA/444/4-89-001, 1989
Figure 6.1-2 pg. 6-3

Benthic Macroinvertebrate Organism Sampling

MI - 4
SAMPLE STATION ID.

MAY 20, 1992
DATE

Rapid Bioassessment Protocol I

Biosurvey Field Data Sheet

RELATIVE ABUNDANCE OF AQUATIC BIOTA

Periphyton	0	1	2	3	4		Slimes	0	1	2	3	4
Filamentous Algae	0	1	2	3	4		Macroinvertebrates	0	1	2	3	4
Macrophytes	0	1	2	3	4		Fish	0	1	2	3	4

0 = Absent/Not Observed 1 = Rare 2 = Common 3 = Abundant 4 = Dominant

MACROBENTHOS QUALITATIVE SAMPLE LIST (indicate Relative Abundance R = Rare, C = Common, A = Abundant, D = Dominant)

Portera	Anisoptera	Chironomidae
Hydrozoa	Zygoptera	Plecoptera D
Platyhelminthes	Hemiptera	Ephemeroptera D
Turbellaria	Coleoptera R	Trichoptera D
Hirudinea	Lepidoptera	Other
Oligochaeta	Sialidae	newt: <i>Desmognathus</i>
Isopoda	Corydalidae	
Amphipoda	Tipulidae R	
Decapoda C	Empididae	
Gastropoda	Simuliidae	
Bivalvia	Tabanidae	
	Culicidae	

Rare < 3 Common 3-8 Abundant > 10 Dominant > 50 (Estimate)

Observations

plant species observed:

christmas fern
honeysuckle
mayapple
sugar maple
tulip tree
bdy's slipper fern

jack in the pulpit
shagbark hickory
sippy elm
rubus sp.

bloodroot

GOLDER PROJECT NO. 913-6773

Reference: Figure taken from USEPA Guidance
Manual No. EPA/444/4-89-001, 1989
Figure 6.1-1 pg. 6-2

Benthic Macroinvertebrate Organism Sampling

MI - 4
SAMPLE STATION ID.

MAY 20 1992
DATE

IMPAIRMENT ASSESSMENT SHEET

1. Detection of impairment: Impairment detected (Complete items 2-6) No impairment detected (Stop here)

2. Biological impairment indicator:

- | | |
|---|--------------------------------------|
| Benthic macroinvertebrates | Other aquatic communities |
| <input type="checkbox"/> absence of EPT taxa | <input type="checkbox"/> Periphyton |
| <input type="checkbox"/> dominance of tolerant groups | <input type="checkbox"/> filamentous |
| <input type="checkbox"/> low benthic abundance | <input type="checkbox"/> other |
| <input type="checkbox"/> low taxa richness | <input type="checkbox"/> Macrophytes |
| <input type="checkbox"/> other | <input type="checkbox"/> Slimes |
| | <input type="checkbox"/> Fish |

3. Brief description of problem: _____
Year and date of previous surveys: _____
Survey data available in: _____

4. Causes (indicate major cause) organic enrichment toxicants flow
habitat limitations other _____

5. Estimated areal extent of problem (m²) and length of stream reach affected (m), where applicable: _____

6. Suspected source(s) of problem:

- point source discharge (name, type of facility, location)
- construction site runoff
- combined sewer outfall
- silviculture runoff
- animal feedlot
- agricultural runoff
- urban runoff
- ground water
- other
- unknown

Briefly explain:

GOLDER PROJECT NO. 913-6773

Reference: Figure taken from USEPA Guidance
Manual No. EPA/444/4-89-001, 1989
Figure 6.1-2 pg. 6-3

Benthic Macroinvertebrate Organism Sampling

M1-5
SAMPLE STATION ID.

MAY 20, 1992
DATE

Rapid Bioassessment Protocol I

Biosurvey Field Data Sheet

RELATIVE ABUNDANCE OF AQUATIC BIOTA

Periphyton	0	1	2	3	4		0	1	2	3	4
Filamentous Algae	0	①	2	3	4	Slimes	0	1	2	3	4
Macrophytes	0	1	2	3	4	Macroinvertebrates	0	1	2	3	④
						Fish	0	1	2	③	4

0 = Absent/Not Observed 1 = Rare 2 = Common 3 = Abundant 4 = Dominant

MACROBENTHOS QUALITATIVE SAMPLE LIST (Indicate Relative Abundance R = Rare, C = Common, A = Abundant, D = Dominant)

Portiera	Anisoptera	Chironomidae	R
Hydrozoa	Zygoptera	Plecoptera	D
Platyhelminthes	Hemiptera	Ephemeroptera	D
Turbellaria	Coleoptera	Trichoptera	R
Hirudinea	Lepidoptera	Other	
Oligochaeta	Sialidae	Newt: <i>Dicamptylus</i> sp.	
Isopoda	Corydalidae		
Amphipoda	Tipulidae		R
Decapoda	Empididae		R
Gastropoda	Simuliidae		R
Bivalvia	Tabanidae		
	Culicidae		

Rare < 3 Common 3-9 Abundant > 10 Dominant > 60 (Estimate)

Observations

plant species observed:

skunk cabbage
christmas fern
honeysuckle
may apple
sugar maple
tulip tree
cedar

garlic mustard
cleaverweed
jack in the pulpit
poison ivy
common burdock

multiflora rose
common reed
larch
curly dock

GOLDER PROJECT NO. 913-6773

Reference: Figure taken from USEPA Guidance
Manual No. EPA/444/4-89-001, 1989
Figure 6.1-1 pg. 6-2

Benthic Macroinvertebrate Organism Sampling

MI - 5
SAMPLE STATION ID.

MAY 20, 1992
DATE

IMPAIRMENT ASSESSMENT SHEET

1. Detection of impairment: Impairment detected (Complete items 2-6) No impairment detected (Stop here)
2. Biological impairment indicator:
- | | |
|---|--------------------------------------|
| Benthic macroinvertebrates | Other aquatic communities |
| <input type="checkbox"/> absence of EPT taxa | <input type="checkbox"/> Periphyton |
| <input type="checkbox"/> dominance of tolerant groups | <input type="checkbox"/> filamentous |
| <input type="checkbox"/> low benthic abundance | <input type="checkbox"/> other |
| <input type="checkbox"/> low taxa richness | <input type="checkbox"/> Macrophytes |
| <input type="checkbox"/> other | <input type="checkbox"/> Slimes |
| | <input type="checkbox"/> Fish |
3. Brief description of problem: _____
Year and date of previous surveys: _____
Survey data available in: _____
4. Cause(s) (indicate major cause) organic enrichment toxicants flow habitat limitations other _____
5. Estimated areal extent of problem (m²) and length of stream reach affected (m), where applicable: _____
6. Suspected source(s) of problem:
- point source discharge (name, type of facility, location)
 - construction site runoff
 - combined sewer outfall
 - silviculture runoff
 - animal feedlot
 - agricultural runoff
 - urban runoff
 - ground water
 - other
 - unknown

Briefly explain:

GOLDER PROJECT NO. 913-6773

Reference: Figure taken from USEPA Guidance
Manual No. EPA/444/4-89-001, 1989
Figure 6.1-2 pg. 6-3

Benthic Macroinvertebrate Organism Sampling

MI-6
SAMPLE STATION ID.

MAY 21 1992
DATE

Rapid Bioassessment Protocol I

Biosurvey Field Data Sheet

RELATIVE ABUNDANCE OF AQUATIC BIOTA

Periphyton	0	1	2	3	4	Slimes	0	1	2	3	4
Filamentous Algae	0	1	2	3	4	Macroinvertebrates	0	1	2	3	4
Macrophytes	0	1	2	3	4	Fish	0	1	2	3	4

0 = Absent/Not Observed 1 = Rare 2 = Common 3 = Abundant 4 = Dominant

MACROBENTHOS QUALITATIVE SAMPLE LIST (Indicate Relative Abundance R = Rare, C = Common, A = Abundant, D = Dominant)

Pterinea	Anisoptera	R	Chironomidae	R
Hydrzoa	Zygoptera		Plecoptera	D
Platyhelminthes	Hemiptera		Ephemeroptera	D
Turbellaria	Coleoptera	C	Trichoptera	R
Miracidae	Lepidoptera		Other	
Oligochaeta	Sialidae			
Isopoda	Corydellidae			
Amphipoda	Tipulidae	R		
Decapoda	Empididae			
Gastropoda	Simuliidae	R		
Bivalvia	Tabanidae			
	Culicidae			

Rare < 3 Common 3-8 Abundant > 10 Dominant > 50 (Estimate)

Observations

Plant species observed:

curly dock
cleared
honeysuckle
ground ivy
garlic mustard
Syramore

ash
multiflora rose
grasses
wild onion
skunk cabbage

Jack in the pulpit
wild geranium
rubus sp.

GOLDER PROJECT NO. 913-6773

Reference: Figure taken from USEPA Guidance
Manual No. EPA/444/4-89-001, 1989
Figure 6.1-1 pg. 6-2

Benthic Macroinvertebrate Organism Sampling

MI-6
SAMPLE STATION ID.

MAY 21 1992
DATE

IMPAIRMENT ASSESSMENT SHEET

1. Detection of impairment: Impairment detected
(Complete items 2-6)

No impairment detected
(Stop here)

2. Biological impairment indicator:

Benthic macroinvertebrates

- absence of EPT taxa
 dominance of tolerant groups
 low benthic abundance
 low taxa richness
 other

Other aquatic communities

- Periphyton
 filamentous
 other
 Macrophytes
 Slimes
 Fish

3. Brief description of problem: _____
Year and date of previous surveys: _____
Survey data available in: _____

4. Cause: (indicate major cause) organic enrichment toxicants flow
habitat limitations other _____

5. Estimated areal extent of problem (m²) and length of stream reach
affected (m), where applicable: _____

6. Suspected source(s) of problem:

- point source discharge (name, type of facility, location)
 construction site runoff
 combined sewer outfall
 silviculture runoff
 animal feedlot
 agricultural runoff
 urban runoff
 ground water
 other
 unknown

Briefly explain:

GOLDER PROJECT NO. 913-6773

Reference: Figure taken from USEPA Guidance
Manual No. EPA/444/4-89-001, 1989
Figure 6.1-2 pg. 6-3

Attachment N-2

Grain Size Distribution Data

AR302682

ASTM GRAIN SIZE ANALYSIS
ASTM D421, D422, D1140, D2216 and D2217

PROJECT TITLE:	BERKS PRP/RIFS/SPRING TWP PA	Sample No:	GSD 1
PROJECT NUMBER:	913-6773	Sample ID:	BULK BAG

WATER CONTENT (Delivered Moisture)		% PASSING #10 SIEVE	
Tare no.	OOP1	Total Wt	2365.84
Wt soil & tare, moist	341.07	Wt Split #10	829.19
Wt soil & tare, dry	318.07	% PASSING #10	22.4%
Wt tare	111.51		
Wt moisture	23.00		
Wt dry soil	206.56		
% WATER	11.13%		

USCS	SIEVE	wt ret	% ret	% PASS	SIEVE
coarse gravel	3.000		0.00%	100.00%	3.000 coarse gravel
	1.500	366.31	15.48%	84.52%	1.500
	1.000	803.43	33.96%	66.04%	1.000
fine gravel	0.750	1019.46	43.09%	56.91%	0.750 fine gravel
	0.375	1377.80	58.24%	41.76%	0.375
coarse sand	#4	1622.40	68.58%	31.42%	#4 coarse sand
medium sand	#10	1836.32	77.62%	22.39%	#10 medium sand

SAMPLE PREPARATION FOR HYDROMETER ANALYSIS	
% Pass #10 Sieve	22.38
Specific Gravity	2.65
ml Dispersing Agent Used	125
Initial Moist Wt.	70.29
Calculated Dry Wt	69.49
(40ml Na(PO4)n per 1000ml H2O)	

WATER CONTENT (Hygroscopic #10)		% PASSING 200 SIEVE	
Tare no.	49	TOTAL WT (g)	
Wt soil and tare, moist	47.62	Wt soil and tare, dry (wt, dry)	188.99
Wt soil & tare, dry	47.32	Wt soil & tare, wash (wt, wash)	183.97
Wt tare	21.13	Wt tare	119.30
Wt moisture lost	0.30	Wt fines lost (wt, dry-wt, wash)	8.02
Wt dry soil, final	26.19	Wt dry soil, (wt, dry-wt, tare)	69.49
% HYGROSCOPIC MOISTURE	1.1%	% FINES LOST	7.2%

PERCENT BETWEEN #10 AND #200 SIEVE CALCULATION			
SIEVE	CUMUL WT. RETAINED	CUMUL WT. RET. CORR	PERCENT PASSING
#10	0.00	241.00	22.39%
#20	22.04	263.04	15.29%
#40	48.72	289.72	6.69%
#60	58.87	299.87	3.42%
#100	62.69	303.69	2.19%
#200	64.56	305.56	1.59%

DATE	TIME	ET (min)	RDNG R	TEMP T	TEMP. COR K	HYD. RDNG. H	Grain Size Percentages
05/28/92	09:13	2.00	8.0	25.00	0.013	4.00	43.1%
	09:15	4.00	8.0	25.00	0.013	4.00	25.5%
	09:19	8.00	7.5	25.00	0.013	4.00	9.0%
	09:26	15.00	7.0	25.00	0.013	4.00	15.7%
	09:41	30.00	6.5	25.00	0.013	3.50	8.1%
	10:11	60.00	6.0	25.00	0.013	4.00	1.6%
	11:11	120.00	8.0	24.50	0.013	4.00	100.00%
	13:11	240.00	8.0	24.00	0.013	4.00	
05/29/92	17:11	490.00	4.5	24.00	0.013	3.50	
	09:11	1440.0	4.0	22.00	0.013	4.00	

ET (min)	RDNG.C	EFF LTH	K	A	PAR DIA	%FINER	WET COLOR: Dark grayish brown
2.00	4.00	15.6	0.013	1.00	0.036	1.3	DESCRIPTION: of GRAVEL, sand & silt. <input type="checkbox"/>
4.00	4.00	15.6	0.013	1.00	0.025	1.3	
8.00	3.50	15.8	0.013	1.00	0.018	1.1	
15.00	3.00	15.8	0.013	1.00	0.013	1.0	
30.00	3.00	15.8	0.013	1.00	0.009	1.0	
60.00	2.00	16.0	0.013	1.00	0.007	0.6	
120.00	1.00	16.1	0.013	1.00	0.005	0.3	
240.00	1.00	16.1	0.013	1.00	0.003	0.3	
490.00	1.00	16.1	0.013	1.00	0.002	0.3	
1440.00	0.00	16.3	0.013	1.00	0.001	0.0	

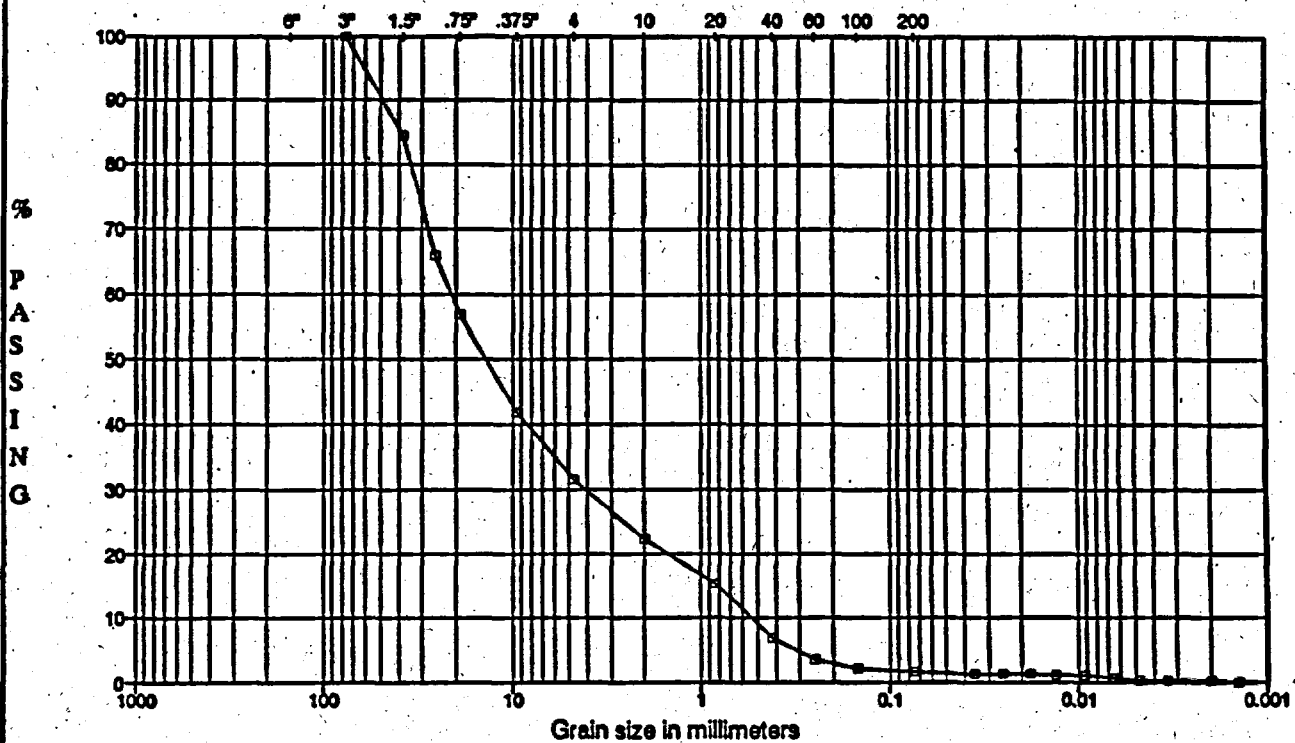
GOLDER ASSOCIATES INC
 MT. LAUREL, NJ

TECH TK
 DATE 5/29/92

CHECKED [Signature]
 REVIEWED [Signature]

AR302683

**PARTICLE SIZE DISTRIBUTION ASTM D421 AND D422
US STANDARD SIEVE OPENING SIZES**



USCS	Coarse	Fine	Co	Med	Fine	FINES (Silt and Clay)
	GRAVEL		SAND			
COBBLES						

TECH: TK
DATE: 6/1/92
CHECKED: *pmu*
REVIEWED:

SAMPLE ID	W%	LL	PL	PI	Gs	DESCRIPTION
GSD 1	11.1					Dark grayish brown c-f GRAVEL, some m sand, trace fines.
Sample Type:	BULK		Date Tested:	5/29/92		

BERKS PRP/RIFS/SPRING TWP PA
913-6773

GOLDER ASSOCIATES INC.
MT. LAUREL, NJ

AR302684

ASTM GRAIN SIZE ANALYSIS
ASTM D421, D422, D1146, D2216 and D2217

PROJECT TITLE:	BERKS PRP/RIFS/SPRING TWP PA	Sample No.:	GSD 2
PROJECT NUMBER:	913-6773	Sample ID.:	BULK BAG

WATER CONTENT (Delivered Moisture)		% PASSING #10 SIEVE	
Tare no.	E-3	Total Wt	1843.84
Wt soil & tare, moist	863.51	Wt Spilt #10	873.77
Wt soil & tare, dry	497.67	% PASSING #10	31.1%
Wt tare	191.53		
Wt moisture	65.84		
Wt dry soil	306.14		
% WATER	21.51%		

U S C S	SIEVE	wt ret	% ret	% PASS	SIEVE
coarse gravel	3.000		0.00%	100.00%	3.000 coarse gravel
	1.500		0.00%	100.00%	1.500
	1.000	179.77	9.75%	90.25%	1.000
fine gravel	0.750	247.84	13.44%	86.56%	0.750 fine gravel
	0.375	649.88	35.25%	64.75%	0.375
coarse sand	#4	1805.13	84.51%	45.49%	#4 coarse sand
medium sand	#10	1270.85	68.92%	31.08%	#10 medium sand

SAMPLE PREPARATION FOR HYDRONETER ANALYSIS			
% Pass #10 Sieve	31.08	Initial Moist Wt.	69.65
Specific Gravity	2.65	Calculated Dry Wt	69.11
ml Dispersing Agent Used	125	(40ml Na(PO4)n per 1800ml H2O)	

WATER CONTENT (Hygroscopic #10)		% PASSING #200 SIEVE	
Tare no.	4B	TOTAL WT (g)	M-2
Wt soil and tare, moist	45.20	Wt soil and tare, dry (wt, dry)	228.68
Wt soil & tare, dry	45.01	Wt soil & tare, wash (wt, wash)	209.68
Wt tare	20.76	Wt tare	151.37
Wt moisture lost	0.19	Wt fines lost (wt, dry-wt, wash)	10.80
Wt dry soil, final	24.25	Wt dry soil, (wt, dry-wt, tare)	69.11
% HYGROSCOPIC MOISTURE	0.9%	% FINES LOST	15.6%

PERCENT BETWEEN #10 AND #200 SIEVE CALCULATION			
SIEVE	CUMUL. WT. RETAINED	CUMUL. WT. RET. CORR	PERCENT PASSING
#10	0.00	153.28	31.08%
#20	14.26	167.54	34.66%
#40	32.01	185.29	36.68%
#60	46.43	199.71	38.20%
#100	54.35	207.63	40.64%
#200	58.14	211.42	43.93%

DATE	TIME	ET (min)	RDNG R	TEMP T	TEMP. COR K	HYDR. RDNG. H	Grain Size Percentages
05/28/92	09:22	2.00	11.0	25.00	0.013	4.00	0.075 13.4%
	09:24	4.00	9.5	25.00	0.013	4.00	0.150 41.1%
	09:28	8.00	9.8	25.00	0.013	4.00	0.300 14.4%
	09:35	15.00	8.5	25.00	0.013	3.50	0.450 14.4%
	09:50	30.00	8.0	25.00	0.013	3.50	0.600 11.7%
	10:20	60.00	7.5	25.00	0.013	4.00	0.750 4.9%
	11:20	120.00	6.0	24.50	0.013	4.00	0.900 100.00%
	13:20	240.00	8.0	24.00	0.013	4.00	
	17:20	480.00	8.0	24.00	0.013	3.50	
05/29/92	09:20	1440.00	4.0	22.00	0.013	4.00	

ET (min)	RDNG. C	EFF LTH	K	A	PAR DIA	% FINER	WET COLOR DESCRIPTION
2.00	7.00	15.2	0.013	1.00	0.035	3.1	Dark grayish brown to black, silty sand and clay. Non-plastic
4.00	5.50	15.5	0.013	1.00	0.025	2.5	
8.00	5.00	15.5	0.013	1.00	0.018	2.2	
15.00	5.00	15.5	0.013	1.00	0.013	2.2	
30.00	4.50	15.6	0.013	1.00	0.009	2.0	
60.00	3.50	15.8	0.013	1.00	0.007	1.6	
120.00	2.00	16.0	0.013	1.00	0.005	0.9	
240.00	1.00	16.1	0.013	1.00	0.003	0.4	
480.00	1.50	16.1	0.013	1.00	0.002	0.7	
1440.00	0.00	16.3	0.013	1.00	0.001	0.0	

GOLDER ASSOCIATES INC MT. LAUREL, NJ	TECH DATE 6/1/92	CHECKED S/W	REVIEWED S/W
---	---------------------	----------------	-----------------

AR302685

ASTM GRAIN SIZE ANALYSIS
ASTM D421, D422, D1144, D2216 and D2217

PROJECT TITLE:	BERKS PRP/RIFS/SPRING TWP PA	Sample No.:	GSD 3
PROJECT NUMBER:	913-6773	Sample ID.:	BULK BAG

WATER CONTENT (Delivered Moisture)		% PASSING #10 SIEVE	
Tare no.	IC-10	Total Wt	1811.52
Wt soil & tare, moist	485.90	Wt Spill #10	293.15
Wt soil & tare, dry	446.11	% PASSING #10	16.2%
Wt tare	190.95		
Wt moisture	45.79		
Wt dry soil	249.16		
% WATER	18.38%		

USCS	SIEVE	wt ret	% ret	% PASS	SIEVE
coarse gravel	3.000		0.00%	100.00%	3.000 coarse gravel
	1.500	94.64	5.22%	94.78%	1.500
	1.000	831.99	29.37%	70.63%	1.000
fine gravel	0.750	828.99	45.76%	54.24%	0.750 fine gravel
	0.375	1212.45	66.93%	33.07%	0.375
coarse sand	#4	1384.73	76.44%	23.56%	#4 coarse sand
medium sand	#10	1518.94	83.85%	16.15%	#10 medium sand

SAMPLE PREPARATION FOR HYDROMETER ANALYSIS			
% Pass #10 Sieve	16.15	Initial Moist Wt	70.80
Specific Gravity	2.65	Calculated Dry Wt	70.12
ml Dispersing Agent Used	125	(40ml Na ₂ PO ₄ in per 1000ml H ₂ O)	

WATER CONTENT (Hygroscopic #10)		% PASSING 200 SIEVE	
Tare no.	1B	TOTAL WT (g)	M-19
Wt soil and tare, moist	86.99	Wt soil and tare, dry (wt, dry)	263.67
Wt soil & tare, dry	86.64	Wt soil & tare, wash (wt, wash)	253.21
Wt tare	28.63	Wt tare	193.85
Wt moisture lost	0.35	Wt fines lost (wt, dry-wt, wash)	8.46
Wt dry soil, final	36.01	Wt dry soil, (wt, dry-wt, tare)	70.12
% HYGROSCOPIC MOISTURE	1.0%	% FINES LOST	12.1%

PERCENT BETWEEN #10 AND #200 SIEVE CALCULATION			
SIEVE	CUMUL. WT. RETAINED	CUMUL. WT. RET. CORR	PERCENT PASSING
#10	0.00	364.02	16.15%
#20	18.65	382.67	11.86%
#40	40.97	404.99	6.71%
#60	83.06	417.08	3.93%
#100	88.64	422.66	2.64%
#200	61.63	425.65	1.96%

DATE	TIME	ET (min)	RDNG R	TEMP T	TEMP. COR K	HYD. RDNG. H	Grain Size Percentages
05/23/92	09:29	2.00	11.0	25.00	0.013	4.00	ccmv. 45.8%
	09:31	4.00	9.5	25.00	0.013	4.00	ccmv. 30.7%
	09:35	8.00	9.0	25.00	0.013	3.50	ccmv. 7.4%
	09:42	15.00	9.0	25.00	0.013	3.50	ccmv. 9.4%
	09:57	36.00	8.0	25.00	0.013	3.50	ccmv. 4.9%
	10:27	68.00	7.5	25.00	0.013	4.00	ccmv. 2.0%
	11:27	128.00	6.0	24.50	0.013	4.00	ccmv. 100.00%
	13:27	240.00	6.5	24.00	0.013	4.00	
	17:27	490.00	6.0	24.00	0.013	4.00	
05/23/92	09:27	1440.0	4.0	22.00	0.013	4.00	

ET (min)	RDNG. C	EFF LTH	K	A	PAR DIA	% FINER	WET COLOR DESCRIPTION
2.00	7.00	15.2	0.013	1.00	0.035	1.6	Dark grayish brown of GRAVEL, some med. trace fines
4.00	6.50	15.5	0.013	1.00	0.025	1.3	
8.00	6.50	15.5	0.013	1.00	0.018	1.3	
15.00	6.50	15.5	0.013	1.00	0.013	1.3	
36.00	4.50	15.6	0.013	1.00	0.009	1.0	
68.00	3.50	15.8	0.013	1.00	0.007	0.8	
128.00	2.00	16.0	0.013	1.00	0.005	0.5	
240.00	1.50	16.1	0.013	1.00	0.003	0.3	
490.00	1.00	16.1	0.013	1.00	0.002	0.2	
1440.00	0.00	16.3	0.013	1.00	0.001	0.0	

GOLDER ASSOCIATES INC
 MT. LAUREL, NJ
 TECH DATE 5/23/92
 CHECKED REVIEWED

AR302687

ASTM GRAIN SIZE ANALYSIS
ASTM D421, D422, D1146, D2216 and D2217

PROJECT TITLE:	BERKS PRP/RIFS/SPRING TWP PA	Sample No.:	GSD 4
PROJECT NUMBER:	913-6773	Sample ID.:	BULK BAG

WATER CONTENT (Delivered Moisture)		% PASSING #10 SIEVE	
Tare no.	2	Total Wt	1353.70
Wt soil & tare, moist	316.78	Wt Split #10	354.93
Wt soil & tare, dry	290.04	% PASSING #10	26.2%
Wt tare	110.85		
Wt moisture	26.74		
Wt dry soil	169.19		
% WATER	21.72%		

U S C S	SIEVE	wt ret	% ret	% PASS	SIEVE
			<small>(% Retention)</small>		
coarse gravel	3.000		0.00%	100.00%	3.000 coarse gravel
	1.500		0.00%	100.00%	1.500
	1.000	65.37	4.83%	95.17%	1.000
fine gravel	0.750	110.26	8.19%	91.85%	0.750 fine gravel
	0.375	495.61	36.61%	63.39%	0.375
coarse sand	#4	781.19	57.71%	42.29%	#4 coarse sand
medium sand	#10	999.41	73.83%	26.17%	#10 medium sand

SAMPLE PREPARATION FOR HYDROMETER ANALYSIS			
% Pass #10 Sieve	26.17	Initial Moist Wt	69.01
Specific Gravity	2.65	Calculated Dry Wt	68.22
ml Dispersing Agent Used	125	(40ml Na(PO4)n per 1000ml H2O)	

WATER CONTENT (Hygroscopic #10)		% PASSING 200 SIEVE	
Tare no.	M-1	TOTAL WT (g)	R-4
Wt soil and tare, moist	52.60	Wt soil and tare, dry (wt, dry)	254.17
Wt soil & tare, dry	52.25	Wt soil & tare, wash (wt, wash)	249.60
Wt tare	22.07	Wt tare	185.95
Wt moisture lost	0.35	Wt fines lost (wt, dry-wt, wash)	4.57
Wt dry soil, final	30.18	Wt dry soil, (wt, dry-wt, tare)	68.22
% HYGROSCOPIC MOISTURE	1.2%	% FINES LOST	6.7%

PERCENT BETWEEN #10 AND #200 SIEVE CALCULATION					
SIEVE	CUMUL. WT. RETAINED	CUMUL. WT. RET. CORR	PERCENT PASSING	LL	PL
#10	0.00	192.44	26.17%		
#20	30.17	222.61	14.60%		
#40	53.52	245.96	5.64%		
#60	60.82	253.26	2.84%		
#100	62.77	255.21	2.09%		
#200	63.75	256.19	1.71%		

DATE	TIME	ET (min)	RDNG R	TEMP T	TEMP. COR K	HYD. RDNG. H	Grain Size Percentages
05/28/92	09:39	2.00	7.0	25.00	0.013	3.50	0.00%
	09:41	4.00	6.5	25.00	0.013	3.50	0.00%
	09:45	8.00	6.5	25.00	0.013	3.50	0.00%
	09:52	15.00	6.5	25.00	0.013	3.50	0.00%
	10:07	30.00	6.0	25.00	0.013	4.00	0.00%
	10:37	60.00	6.0	25.00	0.013	4.00	0.00%
	11:37	120.00	5.0	24.50	0.013	4.00	0.00%
	13:37	240.00	4.0	24.00	0.013	4.00	0.00%
	17:37	490.00	4.0	24.00	0.013	4.00	0.00%
05/29/92	09:37	1440.0	4.0	24.00	0.013	4.00	0.00%

ET (min)	RDNG. C	EFF LTH	K	A	PAR DIA	% FINER	WET COLOR: DESCRIPTION
2.00	3.50	15.8	0.013	1.00	0.026	1.3	Dark grayish brown GRAVEL and SAND. No fines
4.00	3.00	15.8	0.013	1.00	0.026	1.2	
8.00	3.00	15.8	0.013	1.00	0.018	1.2	
15.00	3.00	15.8	0.013	1.00	0.013	1.2	
30.00	2.00	16.0	0.013	1.00	0.009	0.8	
60.00	2.00	16.0	0.013	1.00	0.007	0.8	
120.00	1.00	16.1	0.013	1.00	0.005	0.4	
240.00	0.00	16.3	0.013	1.00	0.003	0.0	
490.00	0.00	16.3	0.013	1.00	0.002	0.0	
1440.00	0.00	16.3	0.013	1.00	0.001	0.0	

GOLDER ASSOCIATES INC
 MT. LAUREL, NJ

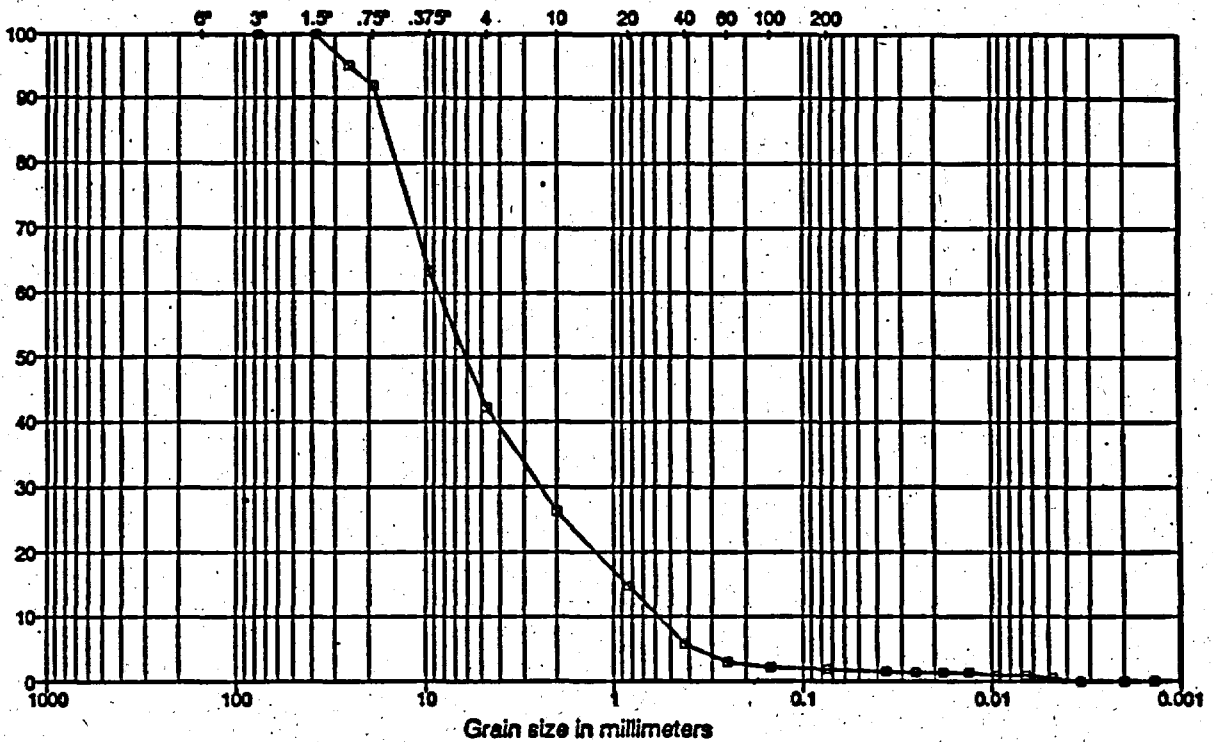
TECH: TK
 DATE: 5/28/92

CHECKED: [Signature]
 REVIEWED: [Signature]

AR302689

**PARTICLE SIZE DISTRIBUTION ASTM D421 AND D422
US STANDARD SIEVE OPENING SIZES**

**P
A
S
S
I
N
G**



USCS

COBBLES	Coarse	Fine	Co	Med	Fine	FINES (Silt and Clay)
	GRAVEL		SAND			

TECH: TK
DATE: 6/1/92
CHECKED: *[Signature]*
REVIEWED:

SAMPLE ID	W%	LL	PL	PI	Gs	DESCRIPTION
GSD 4	21.7					Dark grayish brown f GRAVEL and m-e SAND, trace fines
Sample Type: BULK		Date Tested: 5/28/92				

BERKS PRP/RIFS/SPRING TWP PA
913-6773

GOLDER ASSOCIATES INC.
MT. LAUREL, NJ

AR302690

ASTM GRAIN SIZE ANALYSIS
ASTM D421, D422, D1144, D2216 and D2217

PROJECT TITLE: **BERKS PRP/RIFS/SPRING TWP PA** Sample No: **GSD 5**
 PROJECT NUMBER: **913-6773** Sample ID: **BULK BAG**

WATER CONTENT (Delivered Moisture)			% PASSING #10 SIEVE		
Tare no.		C-40	Total Wt	(g)	1239.91
Wt soil & tare, moist	(g)	582.02	Wt Split #10	(g)	430.83
Wt soil & tare, dry	(g)	512.40	% PASSING #10	(%)	34.7%
Wt tare	(g)	187.35			
Wt moisture	(g)	69.62			
Wt dry soil	(g)	325.05			
% WATER	(%)	21.42%			

USCS	SIEVE	wt ret	% ret	% PASS	SIEVE
coarse gravel	3.000		0.00%	100.00%	3.000 coarse gravel
	1.500		0.00%	100.00%	1.500
	1.000	83.77	6.76%	93.24%	1.000
fine gravel	0.750	149.28	12.04%	87.96%	0.750 fine gravel
	0.375	411.41	33.18%	66.82%	0.375
coarse sand	#4	597.23	48.17%	51.83%	#4 coarse sand
medium sand	#10	810.56	65.37%	34.63%	#10 medium sand

SAMPLE PREPARATION FOR HYDROMETER ANALYSIS

% Pass #10 Sieve	34.63	Initial Moist Wt.	69.55
Specific Gravity	2.65	Calculated Dry Wt	67.97
ml Dispersing Agent Used	125	(40ml Na(PO4)n per 1000ml H2O)	

WATER CONTENT (Hygroscopic #10)			% PASSING 200 SIEVE		
Tare no.		16B	TOTAL WT (g)	(g)	5M3
Wt soil and tare, moist	(g)	43.75	Wt soil and tare, dry (wt, dry)	(g)	262.71
Wt soil & tare, dry	(g)	43.21	Wt soil & tare, wash (wt, wash)	(g)	258.98
Wt tare	(g)	20.02	Wt tare	(g)	194.74
Wt moisture lost	(g)	0.54	Wt fines lost (wt, dry-wt, wash)	(g)	3.73
Wt dry soil, final	(g)	23.19	Wt dry soil, (wt, dry-wt, tare)	(g)	67.97
% HYGROSCOPIC MOISTURE	(%)	2.3%	% FINES LOST	(%)	5.5%

PERCENT BETWEEN #10 AND #200 SIEVE CALCULATION			
SIEVE	CUMUL. WT. RETAINED	CUMUL. WT. RET. CORR	PERCENT PASSING
#10	0.00	128.31	34.63%
#20	40.30	168.61	14.10%
#40	57.23	185.54	8.47%
#60	61.44	189.75	3.33%
#100	63.06	191.37	2.50%
#200	63.97	192.28	2.04%

DATE	TIME	ET (min)	RDNG R	TEMP T	TEMP. COR K	HYD. RDNG. H	Grain Size Percentages
05/28/92	09:48	2.00	7.5	25.00	0.013	3.50	coarse 12.0%
	09:50	4.00	7.5	25.00	0.013	3.50	open 36.1%
	09:54	8.00	7.5	25.00	0.013	3.50	closed 17.3%
	10:01	15.00	7.5	25.00	0.013	3.50	open 29.2%
	10:16	30.00	7.0	25.00	0.013	4.00	open 3.4%
	10:46	60.00	6.0	25.00	0.013	4.00	open 2.0%
	11:46	120.00	5.0	24.50	0.013	4.00	total 100.00%
	13:46	240.00	4.5	24.00	0.013	4.00	
05/29/92	09:46	490.00	4.5	24.00	0.013	3.50	
	09:46	1440.00	4.0	22.00	0.013	4.00	

ET (min)	RDNG. C	EFF LTH	K	A	PAR DIA	% FINER	WET COLOR: DESCRIPTION
2.00	4.00	15.6	0.013	1.00	0.036	2.0	Dark grayish brown non-plastic fine sand
4.00	4.00	15.6	0.013	1.00	0.025	2.0	
8.00	4.00	15.6	0.013	1.00	0.018	2.0	
15.00	4.00	15.6	0.013	1.00	0.013	2.0	
30.00	3.00	15.6	0.013	1.00	0.009	1.5	
60.00	2.00	16.0	0.013	1.00	0.007	1.0	
120.00	1.00	16.1	0.013	1.00	0.005	0.5	
240.00	0.50	16.3	0.013	1.00	0.003	0.3	
490.00	1.00	16.1	0.013	1.00	0.002	0.2	
1440.00	0.00	16.3	0.013	1.00	0.001	0.0	

GOLDER ASSOCIATES INC
 MT. LAUREL, NJ

TECH: **TK**
 DATE: **5/29/92**

CHECKED: **[Signature]**
 REVIEWED: **[Signature]**

AR302691

ASTM GRAIN SIZE ANALYSIS
ASTM D421, D422, D1144, D2216 and D2217

PROJECT TITLE:	BERKS PRP/RIFS/SPRING TWP PA	Sample No.:	GSD 6
PROJECT NUMBER:	913-6773	Sample ID:	BULK BAG

WATER CONTENT (Delivered Moisture)				% PASSING #10 SIEVE			
Tare no.		A-2		Total Wt	(g)	1266.00	
Wt soil & tare, moist	(g)	830.95		Wt Spill #10	(g)	666.59	
Wt soil & tare, dry	(g)	449.09		% PASSING #10	(%)	82.7%	
Wt tare	(g)	193.63					
Wt moisture	(g)	81.86					
Wt dry soil	(g)	255.46					
% WATER	(%)	32.04%					

USCS	SIEVE	wt ret	% ret	% PASS	SIEVE	
			(% Retention)			
coarse gravel	3.000		0.00%	100.00%	3.000	coarse gravel
	1.500		0.00%	100.00%	1.500	
	1.000	81.42	6.43%	93.57%	1.000	
fine gravel	0.750	109.87	8.69%	91.32%	0.750	fine gravel
	0.375	258.15	20.39%	79.61%	0.375	
coarse sand	#4	439.21	34.69%	65.31%	#4	coarse sand
medium sand	#10	600.00	47.39%	52.61%	#10	medium sand

SAMPLE PREPARATION FOR HYDROMETER ANALYSIS			
% Pass #10 Sieve	52.61	Initial Moist Wt.	69.10
Specific Gravity	2.65	Calculated Dry Wt.	66.94
ml Dispersing Agent Used	125	(40ml Na(PO ₄) ₆ per 1000ml H ₂ O)	

WATER CONTENT (Hygroscopic #10)				% PASSING 200 SIEVE			
Tare no.		M-3		TOTAL WT (g)		M-13	
Wt soil and tare, moist	(g)	49.47		Wt soil and tare, dry (wt, dry)	(g)	260.33	
Wt soil & tare, dry	(g)	48.60		Wt soil & tare, wash (wt, wash)	(g)	252.60	
Wt tare	(g)	21.60		Wt tare	(g)	193.39	
Wt moisture lost	(g)	0.87		Wt fines lost (wt, dry-wt, wash)	(g)	7.73	
Wt dry soil, final	(g)	27.00		Wt dry soil, (wt, dry-wt, tare)	(g)	66.94	
% HYGROSCOPIC MOISTURE	(%)	3.2%		% FINES LOST	(%)	11.5%	

PERCENT BETWEEN #10 AND #200 SIEVE CALCULATION				PERCENT PASSING			LL	PL	PI	Co
SIEVE	CUMUL. WT. RETAINED	CUMUL. WT. RET. CORR		PERCENT PASSING						
#10	0.00	60.31		52.61%						
#20	21.27	81.58		35.89%						
#40	41.95	102.26		19.64%						
#60	51.71	112.02		11.97%						
#100	56.95	117.26		7.85%						
#200	58.99	119.30		6.25%						

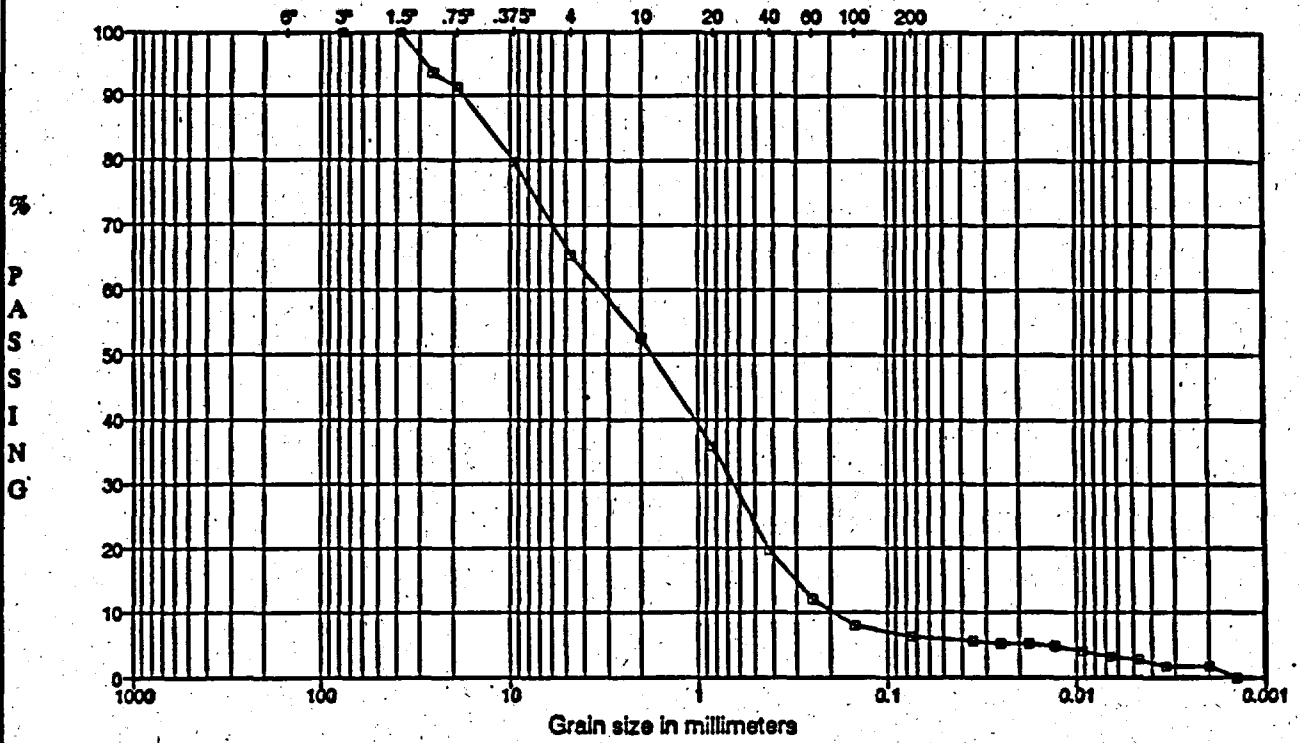
DATE	TIME	ET (min)	RDNG. R	TEMP. T	TEMP. COR. K	HYDRDNG. H	Grain Size Percentages
05/28/92	09:56	2.00	18.5	25.00	0.013	3.50	0.075 8.7%
	09:58	4.00	18.0	25.00	0.013	3.50	0.075 24.0%
	10:02	8.00	18.0	25.00	0.013	3.50	0.075 12.7%
	10:09	15.00	18.0	25.00	0.013	4.00	0.075 33.0%
	10:24	30.00	9.0	25.00	0.013	4.00	0.075 13.4%
	10:54	60.00	8.0	25.00	0.013	4.00	0.075 6.2%
	11:54	120.00	7.5	24.50	0.013	4.00	0.075 100.00%
	13:54	240.00	6.0	24.00	0.013	4.00	
05/29/92	17:54	490.00	8.5	24.00	0.013	3.50	
	09:54	1440.0	4.0	22.00	0.013	4.00	

ET (min)	RDNG. C	EFF LTH	K	A	PAR DIA	% FINER	WET COLOR: DESCRIPTION:
2.00	7.00	15.2	0.013	1.00	0.035	8.5	Dark grayish brown
4.00	6.50	15.3	0.013	1.00	0.025	8.1	Dark grayish brown
8.00	6.50	15.3	0.013	1.00	0.018	8.1	Dark grayish brown
15.00	6.00	15.3	0.013	1.00	0.013	4.7	
30.00	8.00	15.5	0.013	1.00	0.009	3.9	
60.00	4.00	15.6	0.013	1.00	0.007	3.1	
120.00	3.50	15.8	0.013	1.00	0.005	2.8	
240.00	2.00	16.0	0.013	1.00	0.003	1.6	
490.00	2.00	16.0	0.013	1.00	0.002	1.6	
1440.00	0.00	16.3	0.013	1.00	0.001	0.0	

GOLDER ASSOCIATES INC
 MT. LAUREL, NJ
 TECH: TK
 DATE: 5/29/92
 CHECKED: [Signature]
 REVIEWED: [Signature]

AR302693

**PARTICLE SIZE DISTRIBUTION ASTM D421 AND D422
US STANDARD SIEVE OPENING SIZES**



USCS

COBBLES	Coarse	Fine	Co	Med	Fine	FINES (Silt and Clay)
	GRAVEL		SAND			

TECH: TK
DATE: 6/1/92
CHECKED: *[Signature]*
REVIEWED:

SAMPLE ID	W%	LL	PL	PI	Gs	DESCRIPTION
GSD 6	32.0					Dark grayish brown m-f-c SAND and f GRAVEL, little fines
Sample Type: BULK		Date Tested: 5/29/92				

BERKS PRP/RIFS/SPRING TWP PA
913-6773

GOLDER ASSOCIATES INC.
MT. LAUREL, NJ

AR302694

SAMPLE TRACKING RECORD *

JOB NAME: Berks JOB NUMBER: 913-16713 LOCATION: MACKINAC ST. SHEET 1 OF 1
 SAMPLED BY: SAW DATE: 05/21/92 SHIPPED BY: STEPHEN A. WHEELER DATE: MAY 19-21 1992
 SHIP FROM: Berks SHIP TO: GOLDERS SEAL LAB PROJ MRG: SAW / RSN

FIELD		LABORATORY				STORAGE			DISPOSAL			
SAMPLE LOCATION	SAMPLE NUMBER	SAMPLE TYPE (1)	HAZARD (2)	DATE SAMPLE RECEIVED	RECEIVED BY	SAMPLE CONDITION (3)	STORAGE LOCATION	STORAGE DATE	PROJ. MGR. INT.	DISPOSAL DATE	PROJ. MGR. INT.	DISPOSAL METHOD (4)
MI 1	GSD 1	B	*	5/21/92	A. DAVIS		NJ LAB			7/7/92	SAW	R
MI 2	2	B		5/21/92	A. DAVIS		NJ LAB			7/7/92	SAW	R
MI 3	3	B		5/21/92	R. DAVIS		NJ LAB			7/7/92	SAW	R
MI 4	4	B		5/21/92	A. DAVIS		NJ LAB			7/7/92	SAW	R
MI 5	5	B		5/21/92	A. DAVIS		NJ LAB			7/7/92	SAW	R
MI 6	6	B		5/21/92	R. DAVIS		NJ LAB			7/7/92	SAW	R

SPECIAL INSTRUCTION AND NOTES (5)

- o ALL SAMPLES MAY CONTAIN HAZARDOUS CONSTITUENTS
- o ALL SAMPLES TO BE ANALYZED FOR GRAIN SIZE DIST. & 910 MATURE

* THIS FORM IS FOR TRACKING SAMPLES THROUGH THE GOLDERS ASSOCIATES LABORATORY.

ADDITIONAL DOCUMENTATION SHALL BE USED FOR SAMPLES SENT TO OTHER LABORATORIES.

(1) SAMPLE TYPES: J-JAR, T-TUBE, B-BAG, CR-CORE, W-WATER, P-PAL, GX-GEOTEXTILE, GM-GEOMEMBRANE, GM-GEOMET

(2) CODE: C-CHEMICAL, R-RADIOACTIVE, E-EXTREME HAZARD

(3) O-OVER, D-DAMAGED, M-MISSING

(4) T-TRASH, C-CLIENT PICKUP, R-RETURN TO CLIENT, O-OTHER

(5) SPECIAL INSTRUCTIONS SHALL INCLUDE LABORATORY TEST ASSIGNMENTS, PACKING, STORAGE, OR DISPOSAL

#R302696

Attachment N-3

**Laboratory Analytical Results and
Data Validation Forms**

AR302697

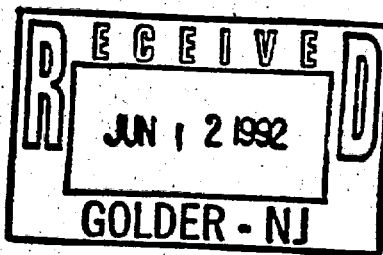
**CEIMIC
CORPORATION**

"Analytical Chemistry for Environmental Management"

June 11, 1992

913 6773.000
(431)

Ms. Lori Hendel
Golder Associates
20000 Horizon Way
Suite 500
Mt. Laurel, NJ 08054



Dear Ms. Hendel:

Enclosed is the data report of results for the analyses of samples received at Ceimic Corporation on May 22, 1992.

Please note that the duplicate and spike analysis for Hardness and Alkalinity were performed after the sample analysis. The analysis date for the Hardness QC set is 6/09/92 and for the Alkalinity QC set it is 6/08/92.

Please don't hesitate to call if you have any questions.

Sincerely,

A handwritten signature in cursive script that reads "Phyllis Shiller".

Phyllis Shiller
Inorganic Laboratory
Manager

PS/11

enc.

**CEIMIC
CORPORATION**

"Analytical Chemistry for Environmental Management"

INORGANIC ANALYTES

Client: Golder Associates

Client ID: SED01/1A

Laboratory ID: 920255-01

Date Sample Received: 5/22/92

Date Sampled: 5/21/92

Target Analyte	Result	Units	Method Reporting Limit	Date Analyzed
Total Organic Carbon (TOC)	0.4	%	0.1	5/29/92

+ Dry weight basis, solid = 76%

Reported by: Account

Approved by: Ayllin Sklar

**CEIMIC
CORPORATION**

"Analytical Chemistry for Environmental Management"

INORGANIC ANALYTES

Client: Golder Associates

Client ID: SED07/1A

Laboratory ID: 920255-03

Date Sample Received: 5/22/92

Date Sampled: 5/20/92

Target Analyte	Result	Units	Method Reporting Limit	Date Analyzed
Total Organic Carbon (TOC)	0.2	%	0.1	5/29/92

+ Dry weight basis, solid = 73%

Reported by: Accident

Approved by: Phyllis Skille

**CEIMIC
CORPORATION**

"Analytical Chemistry for Environmental Management"

INORGANIC ANALYTES

Client: Golder Associates

Client ID: SED08/1A

Laboratory ID: 920255-04

Date Sample Received: 5/22/92

Date Sampled: 5/19/92

Target Analyte	Result	Units	Method Reporting Limit	Date Analyzed
Total Organic Carbon (TOC)	0.2	%*	0.1	5/29/92

+ Dry weight basis, solid = 79%

Reported by: Accountant

Approved by: John Smith

**CEIMIC
CORPORATION**

"Analytical Chemistry for Environmental Management"

INORGANIC ANALYTES

Client: Golder Associates

Client ID: SED09/1A

Laboratory ID: 920255-05

Date Sample Received: 5/22/92

Date Sampled: 5/20/92

Target Analyte	Result	Units	Method Reporting Limit	Date Analyzed
Total Organic Carbon (TOC)	0.3	%⁺	0.1	5/29/92

+ Dry weight basis, solid = 83%

Reported by: Acatest

Approved by: Phyllis Shale

**CEIMIC
CORPORATION**

"Analytical Chemistry for Environmental Management"

QUALITY CONTROL

MATRIX SPIKE ANALYSIS SUMMARY

Client: Golder Associates

Client Sample ID: SED09/1A

Laboratory ID: 920255-05

Date Analyzed: 5/29/92

Concentration in: %

Target Analyte	Sample Result	Spike Added	Spiked Sample Result	Percent Recovery
Total Organic Carbon (TOC)	0.315	0.196	0.506	97%

Reported by: Accutest

Approved by: Byllis Skell

**CEIMIC
CORPORATION**

"Analytical Chemistry for Environmental Management"

QUALITY CONTROL

MATRIX SPIKE DUPLICATE ANALYSIS SUMMARY

Client: Golder Associates

Client Sample ID: SED09/1A

Laboratory ID: 920255-05

Date Analysis Completed: 5/29/92

Target Analyte	Matrix Spike Recovery	Matrix Spike Duplicate Recovery
Total Organic Carbon (TOC)	97%	101%

Reported by: Accutest

Approved by: Phyllis Skell

**CEIMIC
CORPORATION**

"Analytical Chemistry for Environmental Management"

INORGANIC ANALYTES

Client: Golder Associates

Client ID: SED10/1A

Laboratory ID: 920255-06

Date Sample Received: 5/22/92

Date Sampled: 5/19/92

Target Analyte	Result	Units	Method Reporting Limit	Date Analyzed
Total Organic Carbon (TOC)	0.4	%*	0.1	5/29/92

+ Dry weight basis, solid = 74%

Reported by: Accountant

Approved by: Billis Skell

**CEIMIC
CORPORATION**

"Analytical Chemistry for Environmental Management"

INORGANIC ANALYTES

Client: Golder Associates

Client ID: SW-01/1A

Laboratory ID: 920255-07

Date Sample Received: 5/22/92

Date Sampled: 5/21/92

Target Analyte	Result	Units	Method Reporting Limit	Date Analyzed
Alkalinity (as CaCO ₃)	92	mg/L (ppm)	2	5/28/92
Hardness (as CaCO ₃)	151	mg/L (ppm)	2	5/29/92
Total Suspended Solids	13	mg/L (ppm)	5	5/26/92

Reported by:

Jeffrey D. Maxmon

Approved by:

Phyllis Skell

**CEIMIC
CORPORATION**

"Analytical Chemistry for Environmental Management"

INORGANIC ANALYTES

Client: Golder Associates

Client ID: SW-04/1A

Laboratory ID: 920255-08

Date Sample Received: 5/22/92

Date Sampled: 5/20/92

Target Analyte	Result	Units	Method Reporting Limit	Date Analyzed
Alkalinity (as CaCO₃)	72	mg/L (ppm)	2	5/28/92
Hardness (as CaCO₃)	137	mg/L (ppm)	2	5/29/92
Total Suspended Solids	ND	mg/L (ppm)	5	5/26/92

ND = Not detected

Reported by:

J. D. Mayman

Approved by:

Phyllis Skell

CEIMIC
CORPORATION
"Analytical Chemistry for Environmental Management"

INORGANIC ANALYTES

Client: Golder Associates

Client ID: SW-08/1A

Laboratory ID: 920255-09

Date Sample Received: 5/22/92

Date Sampled: 5/19/92

Target Analyte	Result	Units	Method Reporting Limit	Date Analyzed
Alkalinity (as CaCO ₃)	126	mg/L (ppm)	2	5/28/92
Hardness (as CaCO ₃)	188	mg/L (ppm)	2	5/29/92
Total Suspended Solids	ND	mg/L (ppm)	5	5/26/92

ND = Not detected

Reported by:

Jeffrey D. Mayman

Approved by:

Aylin Shale

AR302709

**CEIMIC
CORPORATION**

"Analytical Chemistry for Environmental Management"

INORGANIC ANALYTES

Client: Golder Associates

Client ID: SW-09/1A

Laboratory ID: 920255-10

Date Sample Received: 5/22/92

Date Sampled: 5/19/92

Target Analyte	Result	Units	Method Reporting Limit	Date Analyzed
Alkalinity (as CaCO ₃)	120	mg/L (ppm)	2	5/28/92
Hardness (as CaCO ₃)	163	mg/L (ppm)	2	5/29/92
Total Suspended Solids	23	mg/L (ppm)	5	5/26/92

Reported by:

Jeffrey D. Magnus

Approved by:

Pyllis Skell

**CEIMIC
CORPORATION**

"Analytical Chemistry for Environmental Management"

INORGANIC ANALYTES

Client: Golder Associates

Client ID: SW-10/1A

Laboratory ID: 920255-11

Date Sample Received: 5/22/92

Date Sampled: 5/20/92

Target Analyte	Result	Units	Method Reporting Limit	Date Analyzed
Alkalinity (as CaCO ₃)	92	mg/L (ppm)	2	5/28/92
Hardness (as CaCO ₃)	144	mg/L (ppm)	2	5/29/92
Total Suspended Solids	20	mg/L (ppm)	5	5/26/92

Reported by:

Jeffrey D. Mayman

Approved by:

Phillip Skille

**CEIMIC
CORPORATION**

"Analytical Chemistry for Environmental Management"

QUALITY CONTROL

SPIKE ANALYSIS SUMMARY

Client: Golder Associates

Client Sample ID: SW-10/1A

Laboratory ID: 920255-11

Date Sample Received: 5/22/92

Concentration in: mg/L

Target Analyte	Sample Result	Predigest Spike Added	Spiked Sample Result	Percent Recovery		
				Predigest Spike	Control Limit	Post Digest Spike Recovery*
Alkalinity (as CaCO ₃)	92	25	114	88	75-125%	NR
Hardness (as CaCO ₃)	144	182	361	119	75-125	NR

NR = Not required

* Post digest spike added at twice the CRDL or twice the indigenous level which ever is greater.

Reported by: Jeffrey D. Maymon

Approved by: Phyllis Skell

CEIMIC
CORPORATION
"Analytical Chemistry for Environmental Management"
QUALITY CONTROL

DUPLICATE ANALYSIS SUMMARY

Client: Golder Associates

Client Sample ID: SW-10/1A

Laboratory ID: 920255-11

Date Samples Received: 5/22/92

Date Sampled: 5/20/92

Target Analyte	Sample Result	Duplicate Result	RPD	Control Limit
Alkalinity (as CaCO ₃)	92	91	1%	20%
Hardness (as CaCO ₃)	144	168	15%	20%
Total Suspended Solids	20	18	11%	20%

RPD = Relative Percent Difference

Reported by: *J. D. Mayman*

Approved by: *Byllis Skale*

**CEIMIC
CORPORATION**

"Analytical Chemistry for Environmental Management"

INORGANIC ANALYTES

Client: Golder Associates

Client ID: SW-11/1A

Laboratory ID: 920255-12

Date Sample Received: 5/22/92

Date Sampled: 5/20/92

Target Analyte	Result	Units	Method Reporting Limit	Date Analyzed
Alkalinity (as CaCO₃)	120	mg/L (ppm)	2	5/28/92
Hardness (as CaCO₃)	139	mg/L (ppm)	2	5/29/92
Total Suspended Solids	23	mg/L (ppm)	5	5/26/92

Reported by:

Jeffrey D. Magnuson

Approved by:

Billy Skell

**CEIMIC
CORPORATION**

"Analytical Chemistry for Environmental Management"

INORGANIC ANALYTES

Client: Golder Associates

Client ID: RB001/1A

Laboratory ID: 920255-13

Date Sample Received: 5/22/92

Date Sampled: 5/21/92

Target Analyte	Result	Units	Method Reporting Limit	Date Analyzed
Total Organic Carbon (TOC)	ND	mg/L (ppm)	1	5/28/92

ND = Not detected

Reported by: *Acaster*

Approved by: *Pyllis Skell*

**CEIMIC
CORPORATION**

"Analytical Chemistry for Environmental Management"

QUALITY CONTROL

METHOD BLANK

Client: Golder Associates

Client ID: Method Blank

Project No.: 920255

Laboratory ID: PBS

Concentration in: %

Target Analyte	Result	Method Reporting Limit	Date Analyzed
Total Organic Carbon (TOC)	ND	0.1	5/29/92

ND = Not detected

Reported by: Accutest

Approved by: Phyllis Skell

**CEIMIC
CORPORATION**

"Analytical Chemistry for Environmental Management"

QUALITY CONTROL

LABORATORY CONTROL SAMPLE

Client: Golder Associates

Client Sample ID: Laboratory Control Sample

Project No.: 920255

Laboratory ID: LCSS

Matrix: Soil

Target Analyte	± Recovery	Control Limits
Total Organic Carbon (TOC)	103±	75-125±

Reported by: Accutest

Approved by: Peggy Skelton

**CEIMIC
CORPORATION**

"Analytical Chemistry for Environmental Management"

QUALITY CONTROL

METHOD BLANK

Client: Golder Associates

Client ID: Method Blank

Project No.: 920255

Laboratory ID: PBW

Target Analyte	Result	Units	Method Reporting Limit	Date Analyzed
Alkalinity (as CaCO ₃)	ND	mg/L (ppm)	2	5/28/92
Hardness (as CaCO ₃)	ND	mg/L (ppm)	2	6/10/92
Total Organic Carbon	ND	mg/L (ppm)	1	5/28/92
Total Suspended Solids	ND	mg/L (ppm)	5	5/26/92

ND = Not detected

Reported by:

Jeffrey D. Maymon

Approved by:

Pyllis Stalle

**CEIMIC
CORPORATION**

"Analytical Chemistry for Environmental Management"

QUALITY CONTROL

LABORATORY CONTROL SAMPLE

Client: Golder Associates

Client Sample ID: Laboratory Control Sample

Project No.: 920255

Laboratory ID: LCSW

Matrix: Aqueous

Target Analyte	Recovery	Control Limits
Alkalinity (as CaCO ₃)	101%	75-125%
Hardness (as CaCO ₃)	101	75-125
Total Organic Carbon (TOC)	100	75-125
Total Suspended Solids	84	75-125

Reported by: Jeffrey D. Mayman

Approved by: Pyllis Skell

CHAIN OF CUSTODY

Original chain of Custody goes to Laboratory

Proj. #	Project name	Date	Comp. Grab	Sample Identification	Type of container	Number of containers	Date/Time	Remarks
012-16713	PEAKS DRP RILEY SPRING TOWNSHIP PA							
Samplers (Please print) STEVEN A. WHEELER								
DANIEL J. MILLNER								
							HNO3 NO PRESV.	
5/21 0910	X SN-01/1A			HARDNESS	250 P	1	X	MAXIMUM SBT 10
5/21 0910	SN-01/1A			ALKALINITY TSS	500 P	1	X	10
5/16 1220	SN-04/1A			HARDNESS	250 P	1	X	4
5/20 1220	SN-04/1A			ALKALINITY TSS	500 P	1	X	4
5/19 1015	SN-08/1A			HARDNESS	250 P	1	X	1
5/19 1015	SN-08/1A			ALKALINITY TSS	500 P	1	X	1
5/19 15:00	SN-09/1A			HARDNESS	250 P	1	X	2
5/19 15:00	SN-09/1A			ALKALINITY TSS	500 P	1	X	2
5/20 0840	SN-10/1A			HARDNESS	250 P	1	X	3
5/20 0840	SN-10/1A			ALKALINITY TSS	500 P	1	X	3
5/20 1530	SN-11/1A			HARDNESS	250 P	1	X	5
5/20 1530	SN-11/1A			ALKALINITY TSS	500 P	1	X	5
5/21 0940	R0001/1A			RAINWATER CLEAN	250 P	1		
Refiniquished by (Signature) <i>Daphne A. W. [Signature]</i> Date/Time 05/31/12 Received by (Signature) _____ Date/Time 11/01/15								
Refiniquished by (Signature) _____ Date/Time _____ Received by (Signature) _____ Date/Time _____								
Refiniquished by (Signature) _____ Date/Time _____ Received by (Signature) <i>Ryan R. [Signature]</i> Date/Time 5/22/02								
_____ Date/Time 10/15								

AR302720

P = Plastic

903121

CHAIN OF CUSTODY

Original chain of Custody goes to Laboratory

Project name		Sample Identification		Type of container	Number of containers	Remarks
Date	Time	Comp.	Grab			
5/21	9:20	50		402	1	EH1414B
5/20	12:30				1	EH1414H
5/20					1	EH1414S
5/20					1	EH1414S
5/20	15:10				1	EH1414S
5/19	10:20				1	EH1414B
5/20	8:50				1	EH1414H
5/19	15:10				1	EH1414B
5/21	8:30			3921	1	EH1414H
(17)						

Refriniquished by (Signature)	Date/Time	Received by (Signature)	Date/Time	Remarks:
<i>Rgr. Rham</i>	5/22/92 3:30	Airborne Express	5/22/92 3:30	* Refer to list of Requirements enclosed. * 1st MSD Recoveries & Results LCS + LCS duplicate Recoveries & Results
	5/23/92 12:20		5/23/92 12:20	

Project name: *CEIMIC*

Project name: *CEIMIC*

Date: 5/21/92

Time: 9:20

Comp.: 50

Grab: 01

Sample ID: 920255

Container: 402

Remarks: EH1414B

Remarks: EH1414H

Remarks: EH1414S

Remarks: EH1414S

Remarks: EH1414S

Remarks: EH1414B

Remarks: EH1414H

Remarks: EH1414B

Remarks: EH1414H

Remarks: EH1414H

AR302722

CEIMIC Corporation 10 Dean Knauss Drive, Narragansett, RI 02882 (401) 782-8900 FAX (401) 782-8905

* To Results by 6:10 (1st Day T/A)
* TI-I pg. to follow shortly after.

TI-1

SAMPLES RECEIVED ON ICE AT ACCOUNT

* Results Due 6/10/92

WV ST/AR

Data Evaluation Checklist

Project Name: BEEKS LANDFILL RIFS Project Number: 913-6773.1804
 Reviewer: Lori Anne Hendel Date: July 8, 1992
 Signature: Lori Anne Hendel

Overall Data Quality Assessment

Sample Points: SED011A, SED021A, SED071A, SED081A, SED091A, SED101A
RB0011A, SW-011A, SW-041A, SW-081A, SW-091A, SW-101A, SW-111A

	Y	N	NA
--	---	---	----

1. General
- (a) Were the FSP, QAPP, laboratory reports and field documentation available to support data assessment procedures? ✓ _____ _____
2. Precision (reference FSP and QAPP)
- (a) Are lab duplicate RPD within control limits? _____ _____ _____
- (b) Are field duplicate RPD within control limits? _____ _____ _____
- (c) Are MS/MSD RPD within control limits? _____ _____ ✓
- Overall assessment of precision: Precision was deemed acceptable

3. Accuracy (reference FSP and QAPP)
- (a) Are MS/MSD recoveries within control limits? _____ _____ _____
- (b) Are lab spike (LCS/DCS) recoveries within control limits? _____ _____ _____
- Overall assessment of accuracy: Accuracy of data was deemed acceptable

4. Representativeness
- (a) Were procedures in the FSP followed? If not, were procedural variances approved and documented? _____ _____ _____
- (b) Were sample preservation procedures followed? _____ _____ _____
- (c) Were data reported in the proper units? _____ _____ _____
- (d) Was blank contamination not evident, or well documented at low levels? _____ _____ _____
- (e) Were field duplicates within control limits? _____ _____ ✓
- Overall assessment of representativeness: Representativeness of data was deemed acceptable

AR302723

	Y	N	NA
5. Comparability			
(a) Are data traceable to a standard method?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>
(b) Are methods approved/accepted as producing valid results?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>
(c) Are data reported in proper units?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>
Overall assessment of comparability:	<u>Comparability of data was deemed acceptable</u>		

6. Completeness (reference FSP and QAPP)			
(a) Is the fraction of valid data within control limits?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>
(b) If not, are the data sufficient to meet the task objectives? (If yes, justify below)	<input type="checkbox"/>	<input type="checkbox"/>	<input checked="" type="checkbox"/>
(c) Are critical (background) samples sufficient and have valid results?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>
Overall assessment of completeness:	<u>Completeness of data was deemed acceptable</u>		

7. DQOs (reference FSP and QAPP)			
(a) Are the data usable and consistent with the project data quality objectives?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>

8. Comments/Notes

All data was deemed acceptable based upon the criteria established by the FSP and QAPP

Data Evaluation Checklist

Project Name: BEERS LANDFILL RIFPS Project Number: 913-16713.1004
 Reviewer: Lori Anne Herdel Date: July 8, 1992
 Signature: Lori Anne Herdel

Field Performance

Sample Points: SED011A, SED031A, SED071A, SED081A, SED091A, SED101A
SED011A, SW-011A, SW-041A, SW-081A, SW-091A, SW-101A, SW-111A

	Y	N	NA
1. Chain-of-Custody (CoC)			
(a) Was the CoC properly completed?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>
(b) Was the CoC signed by both field and laboratory personnel?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>
(c) Were samples received in good condition?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>
2. Field Documentation includes (reference FSP)			
(a) Date and times samples collected?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>
(b) Sample locations?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>
(c) Field measurements?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>
(d) Instruments for field measurements (including instrument model and serial number)?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>
(e) Calibration/maintenance of field instruments (including lot number, source and frequency)?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>
(f) Proper preservatives used?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>
(g) Field quality control procedures?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>
(h) FSP procedures followed? If not, procedural variances approved and documented?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>
3. Precision (reference FSP)			
(a) Were Field Duplicate precision criteria met?	<input type="checkbox"/>	<input type="checkbox"/>	<input checked="" type="checkbox"/>
Note: If one or both results were not detected, then RPD can not be calculated and indicate NA.			
4. Blanks			
(a) Contamination of field blanks was not evident, or well documented at low levels?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>

5. Comments/Notes

Data Evaluation Checklist

Project Name: Berks Landfill RIFs Project Number: 913-6773.1004

Reviewer: Lori Anne Hendel Date: June 24, 1992

Signature: Lori Anne Hendel

Inorganic Conventional Parameters

Parameter: Total Organic Carbon Method: Modified EPA 415.1

Matrix: Sediment

Sample Points: SED01/1A, SED03/1A, SED07/1A, SED08/1A, SED09/1A, SED10/1A

	Y	N	NA
1. Chain-of-Custody (CoC)			
(a) Was the CoC properly completed?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>
(b) Was the CoC signed by both field and laboratory personnel?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>
(c) Were samples received in good condition?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>
2. Samples (reference QAPP or Method)			
(a) Were hold times met for all samples?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>
(b) Were the correct preservatives used?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>
(c) Was the correct method used?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>
(d) Were appropriate reporting limits achieved?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>
(e) Were any sample dilutions noted?	<input type="checkbox"/>	<input checked="" type="checkbox"/>	<input type="checkbox"/>
(f) Were any matrix problems noted?	<input type="checkbox"/>	<input checked="" type="checkbox"/>	<input type="checkbox"/>
3. Precision (reference QAPP)			
(a) Were Field Duplicate precision criteria met?	<input type="checkbox"/>	<input type="checkbox"/>	<input checked="" type="checkbox"/>
(b) Were MS/MSD precision criteria met? 4%	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>
(c) Were Lab Duplicate precision criteria met?	<input type="checkbox"/>	<input type="checkbox"/>	<input checked="" type="checkbox"/>

Note: If one or both results were not detected, then RPD can not be calculated and indicate NA.

	Y	N	NA
4. Accuracy (reference QAPP)			
(a) Was MS accuracy criteria met? 97% Recovery could not be calculated since sample contained high concentration of analyte?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>
(b) Was MSD accuracy criteria met? 101% Recovery could not be calculated since sample contained high concentration of analyte?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input checked="" type="checkbox"/>
(c) Was LCS accuracy criteria met? 103% Recovery could not be calculated since sample contained high concentration of analyte?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input checked="" type="checkbox"/>
5. Blanks			
(a) Was analyte detected in the method blank(s)?	<input type="checkbox"/>	<input checked="" type="checkbox"/>	<input type="checkbox"/>
(b) Was analyte detected in the field blank(s)?	<input type="checkbox"/>	<input checked="" type="checkbox"/>	<input type="checkbox"/>

6. Comments/Notes

Pinstate blank collected: GC for aqueous TOC acceptable
(Blank - ND²; LCS 100% recovery)

Samples subcontracted to Accutest

Data Evaluation Checklist

Project Name: Berks Landfill R/FS Project Number: 913-6773.1004
 Reviewer: Lori Anne Hendel Date: June 24, 1992
 Signature: Lori Anne Hendel

Inorganic Conventional Parameters

Parameter: Alkalinity Method: EPA 310.2
 Matrix: Surface Water
 Sample Points: SW-01/1A, SW-04/1A, SW-08/1A, SW-09/1A, SW-10/1A, SW-11/1A

	Y	N	NA
1. Chain-of-Custody (CoC)			
(a) Was the CoC properly completed?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>
(b) Was the CoC signed by both field and laboratory personnel?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>
(c) Were samples received in good condition?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>
2. Samples (reference QAPP or Method)			
(a) Were hold times met for all samples?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>
(b) Were the correct preservatives used?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>
(c) Was the correct method used?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>
(d) Were appropriate reporting limits achieved?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>
(e) Were any sample dilutions noted?	<input type="checkbox"/>	<input checked="" type="checkbox"/>	<input type="checkbox"/>
(f) Were any matrix problems noted?	<input type="checkbox"/>	<input checked="" type="checkbox"/>	<input type="checkbox"/>
3. Precision (reference QAPP)			
(a) Were Field Duplicate precision criteria met?	<input type="checkbox"/>	<input type="checkbox"/>	<input checked="" type="checkbox"/>
(b) Were MS/MSD precision criteria met?	<input type="checkbox"/>	<input type="checkbox"/>	<input checked="" type="checkbox"/>
(c) Were Lab Duplicate precision criteria met? 1%	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>

Note: If one or both results were not detected, then RPD can not be calculated and indicate NA.

	Y	N	NA
4. Accuracy (reference QAPP)			
(a) Was MS accuracy criteria met? 88%	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>
Recovery could not be calculated since sample contained high concentration of analyte?	<input type="checkbox"/>	<input type="checkbox"/>	<input checked="" type="checkbox"/>
(b) Was MSD accuracy criteria met?	<input type="checkbox"/>	<input type="checkbox"/>	<input checked="" type="checkbox"/>
Recovery could not be calculated since sample contained high concentration of analyte?	<input type="checkbox"/>	<input type="checkbox"/>	<input checked="" type="checkbox"/>
(c) Was LCS accuracy criteria met? 101%	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>
Recovery could not be calculated since sample contained high concentration of analyte?	<input type="checkbox"/>	<input type="checkbox"/>	<input checked="" type="checkbox"/>
5. Blanks			
(a) Was analyte detected in the method blank(s)?	<input type="checkbox"/>	<input checked="" type="checkbox"/>	<input type="checkbox"/>
(b) Was analyte detected in the field blank(s)?	<input type="checkbox"/>	<input type="checkbox"/>	<input checked="" type="checkbox"/>

6. Comments/Notes

GC samples analyzed 6/18/92

Data Evaluation Checklist

Project Name: Berks Landfill P/FS Project Number: 913-6773.1004

Reviewer: Lori Anne Hendel Date: June 24, 1992

Signature: Lori Anne Hendel

Inorganic Conventional Parameters

Parameter: Total Suspended Solids Method: EPA 160.1

Matrix: Surface Water

Sample Points: SW-01/1A, SW-04/1A, SW-08/1A, SW-09/1A, SW-10/1A, SW-11/1A

	Y	N	NA
1. Chain-of-Custody (CoC)			
(a) Was the CoC properly completed?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>
(b) Was the CoC signed by both field and laboratory personnel?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>
(c) Were samples received in good condition?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>
2. Samples (reference QAPP or Method)			
(a) Were hold times met for all samples?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>
(b) Were the correct preservatives used?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>
(c) Was the correct method used?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>
(d) Were appropriate reporting limits achieved?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>
(e) Were any sample dilutions noted?	<input type="checkbox"/>	<input checked="" type="checkbox"/>	<input type="checkbox"/>
(f) Were any matrix problems noted?	<input type="checkbox"/>	<input checked="" type="checkbox"/>	<input type="checkbox"/>
3. Precision (reference QAPP)			
(a) Were Field Duplicate precision criteria met?	<input type="checkbox"/>	<input type="checkbox"/>	<input checked="" type="checkbox"/>
(b) Were MS/MSD precision criteria met?	<input type="checkbox"/>	<input type="checkbox"/>	<input checked="" type="checkbox"/>
(c) Were Lab Duplicate precision criteria met? 11%	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>

Note: If one or both results were not detected, then RPD can not be calculated and indicate NA.

Data Evaluation Checklist

Project Name: Berks Landfill E/FS Project Number: 913-6773.1004

Reviewer: Lori Anne Hendel Date: June 24, 1992

Signature: Lori Anne Hendel

Inorganic Conventional Parameters

Parameter: Hardness Method: EPA 200.7

Matrix: Surface Water

Sample Points: SW-01/1A, SW-04/1A, SW-08/1A, SW-09/1A, SW-10/1A, SW-11/1A

	Y	N	NA
1. Chain-of-Custody (CoC)			
(a) Was the CoC properly completed?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>
(b) Was the CoC signed by both field and laboratory personnel?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>
(c) Were samples received in good condition?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>
2. Samples (reference QAPP or Method)			
(a) Were hold times met for all samples?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>
(b) Were the correct preservatives used?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>
(c) Was the correct method used?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>
(d) Were appropriate reporting limits achieved?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>
(e) Were any sample dilutions noted?	<input type="checkbox"/>	<input checked="" type="checkbox"/>	<input type="checkbox"/>
(f) Were any matrix problems noted?	<input type="checkbox"/>	<input checked="" type="checkbox"/>	<input type="checkbox"/>
3. Precision (reference QAPP)			
(a) Were Field Duplicate precision criteria met?	<input type="checkbox"/>	<input type="checkbox"/>	<input checked="" type="checkbox"/>
(b) Were MS/MSD precision criteria met?	<input type="checkbox"/>	<input type="checkbox"/>	<input checked="" type="checkbox"/>
(c) Were Lab Duplicate precision criteria met? 15%	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>

Note: If one or both results were not detected, then RPD can not be calculated and indicate NA.

	Y	N	NA
4. Accuracy (reference QAPP)			
(a) Was MS accuracy criteria met? 119%	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>
Recovery could not be calculated since sample contained high concentration of analyte?	<input type="checkbox"/>	<input type="checkbox"/>	<input checked="" type="checkbox"/>
(b) Was MSD accuracy criteria met?	<input type="checkbox"/>	<input type="checkbox"/>	<input checked="" type="checkbox"/>
Recovery could not be calculated since sample contained high concentration of analyte?	<input type="checkbox"/>	<input type="checkbox"/>	<input checked="" type="checkbox"/>
(c) Was LCS accuracy criteria met? 101%	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>
Recovery could not be calculated since sample contained high concentration of analyte?	<input type="checkbox"/>	<input type="checkbox"/>	<input checked="" type="checkbox"/>
5. Blanks			
(a) Was analyte detected in the method blank(s)?	<input type="checkbox"/>	<input checked="" type="checkbox"/>	<input type="checkbox"/>
(b) Was analyte detected in the field blank(s)?	<input type="checkbox"/>	<input type="checkbox"/>	<input checked="" type="checkbox"/>

6. Comments/Notes

AC samples analyzed 6/9/92

APPENDIX O
AGENCY CORRESPONDENCE FOR NEPA ISSUES

AR302734

Golder Associates Inc.

305 Fellowship Road, Suite 200
Mt. Laurel, NJ USA 08054
Tel: (609) 273-4110
Fax (609) 273-0778



October 27, 1993

Project No.: 913-6773

PA Department of Environmental Resources
Forest Advisory Services/PNDI
P.O. Box 8552
Harrisburg, PA 17105-8552

Attn: Ms. Kathy McKenna

RE: PENNSYLVANIA NATURAL DIVERSITY INVENTORY (PNDI)
SEARCH REQUEST FOR THE BERKS LANDFILL SITE
BERKS COUNTY, PENNSYLVANIA

Ladies and Gentlemen:

Golder Associates Inc. (Golder) has enclosed the Pennsylvania Natural Diversity Inventory (PNDI) search request form for the Berks Landfill Site, Spring Township, Berks County, Pennsylvania. Also enclosed is a Site Location Map showing the Site location on the Sinking Spring, Pennsylvania, 7.5 minute USGS topographic quadrangle map.

Golder requests any information available regarding protected species within a one-mile radius of the Site.

If you should require any additional information, please do not hesitate to contact us.

Very truly yours,

GOLDER ASSOCIATES INC.

A handwritten signature in black ink, appearing to read "Stephen A. Wheeler".

Stephen A. Wheeler
Project Environmental Scientist

A handwritten signature in black ink, appearing to read "Randolph S. White".

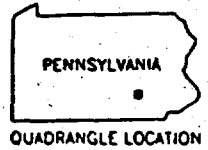
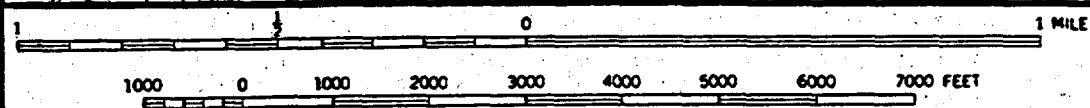
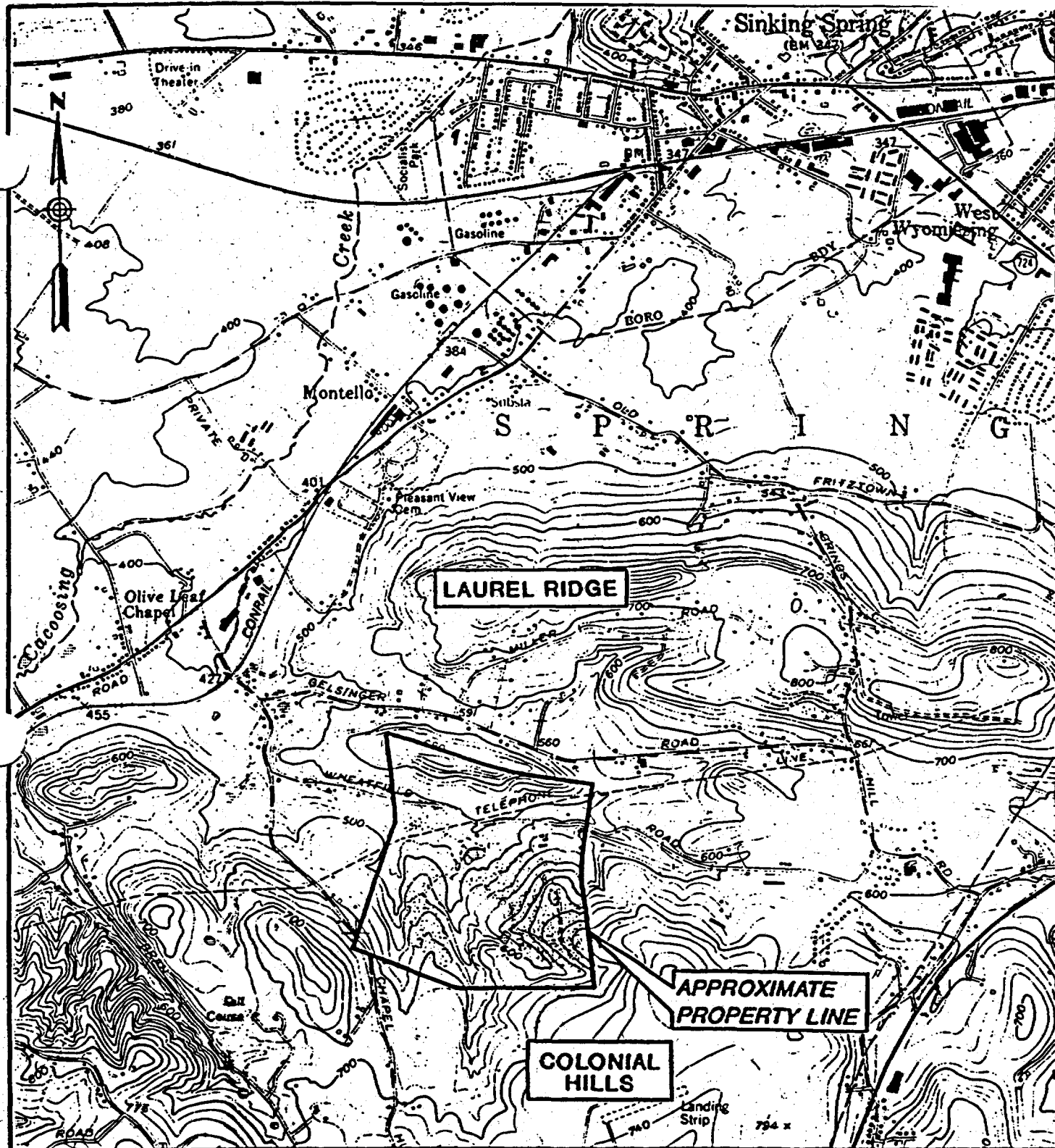
Randolph S. White, P.E.
Senior Project Manager

SAW/RSW:lr1

Enclosure

z:677310271lr3.wp5

AR302735



REFERENCE:
 U.S.G.S. TOPOGRAPHIC MAP, SINKING SPRING QUADRANGLE, 7.5 SERIES, PHOTOREVISED 1987.

DS No.:	913-6389	SCALE:	AS SHOWN
DRAWN:	EAM	DATE:	05/10/91
CHECKED:		DWG. No.:	PA21-075

OCT 27 1993

SITE LOCATION MAP

Golder Associates

BERKS LANDFILL

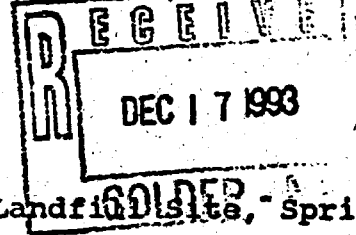
FIGURE

AR302736

COMMONWEALTH OF PENNSYLVANIA
DEPARTMENT OF ENVIRONMENTAL RESOURCES

P.O. Box 8552
Harrisburg, PA 17105-8552

December 14, 1993



717/787-3444



Bureau of Forestry

SUBJECT: PNDI Review of the Berks Landfill Site, Spring Township, Berks County, PA

TO: Stephen A. Wheeler
Golder Associates, Inc.
305 Fellowship Road, Suite 200
Mt. Laurel, NJ 08054

FROM: Pennsylvania Natural Diversity Inventory System

In response to your request of October 27, an area mapped on the Sinking Spring Quadrangle was reviewed for the presence of natural resources of special concern using the Pennsylvania Natural Diversity Inventory (PNDI) information system. We do not anticipate any impact on rare, threatened or endangered species at this location.

PNDI is a site specific information system which describes significant natural resources of Pennsylvania. PNDI includes data descriptive of plant and animal species of special concern, exemplary natural communities and unique geological features. This response represents the most up-to-date summary of the PNDI data files. However, an absence of recorded information does not necessarily imply actual conditions on-site. A field survey of any site may reveal previously unreported populations.

Be advised that legal authority for Pennsylvania's biological resources resides with three administrative agencies. The enclosure titled PNDI Species List, outlines which species groups are managed by these agencies. If you have questions concerning this response or the PNDI system, please contact Edward Dix of our office at 717/783-0392.

Enclosure

cc: Anthony Wilkinson, PNDI-East
File

PENNSYLVANIA NATURAL DIVERSITY INVENTORY

SPECIES LISTS

The statutory authority for Pennsylvania's animals and plants resides with three separate agencies. The Pennsylvania Department of Environmental Resources has the responsibility for management of the Commonwealth's native wild plants. The Pennsylvania Fish Commission is responsible for management of fish, reptiles, amphibians and aquatic organisms within the Commonwealth. The Pennsylvania Game Commission has the responsibility for managing the state's wild birds and mammals.

For information on current official status for a species, please consult the appropriate agency. Requests for information should be directed to:

PLANTS and
PNDI - general

Plant Program Manager
Pa. Department of Environmental Resources
Bureau of Forestry
Forest Advisory Services
P. O. Box 8552
Harrisburg, PA 17105-8552
(717)787-3444

FISH, REPTILES,
AMPHIBIANS,
AQUATIC ORGANISMS

Endangered Species & Herpetology
Coordinator
Pennsylvania Fish & Boat Commission
Bureau of Fisheries and Engineering
450 Robinson Lane
Bellefonte, PA 16823
(814)359-5113

BIRDS and MAMMALS

Pennsylvania Game Commission
Bureau of Wildlife Management
2001 Elmerton Avenue
Harrisburg, PA 17110-9797
(717)787-5529

For information on species listed under the federal Endangered Species Act of 1973 occurring in Pennsylvania, contact:

Endangered Species Biologist
U.S. Fish and Wildlife Service
315 South Allen Street, Suite 322
State College, PA 16801
(814)234-4090

Thank you for your request. Feel free to contact PNDI if we can be of further assistance.

AR302738

Golder Associates Inc.

305 Fellowship Road, Suite 200
Mt. Laurel, NJ USA 08054
Tel: (609) 273-1110
Fax (609) 273-0778



October 27, 1993

Project No.: 913-6773

Pennsylvania Historical and Museum Commission
Bureau for Historic Preservation
William Penn Memorial Museum Building
P.O. Box 1026
Harrisburg, PA 17108-1026

RE: BERKS LANDFILL
SPRING TOWNSHIP, BERKS COUNTY PENNSYLVANIA
HISTORICAL, ARCHEOLOGICAL, AND CULTURAL RESOURCES

Ladies & Gentlemen:

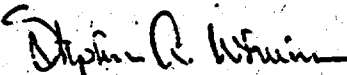
Golder Associates Inc. (Golder) requests a review and determination as to whether any known historical, archeological or cultural resources exist on or within a one-mile radius of the Berks Landfill Site (Site). The Site is located along Wheatfield Road, Spring Township, Berks County, Pennsylvania, as shown on the enclosed Sinking Spring, Pennsylvania 7.5 minute USGS topographic quadrangle map.

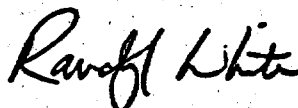
The surface soils at the site have been disturbed from the former construction of a municipal solid waste landfill facility, ancillary facilities and from soil borrow areas utilized by the former landfill. Surrounding areas include farmland, forestland, residential homes, roads and other developments.

If you should require additional information for this request, please do not hesitate to contact us at (609) 273-1110.

Very truly yours,

GOLDER ASSOCIATES INC.


Stephen A. Wheeler
Project Environmental Scientist

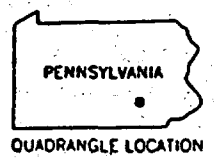
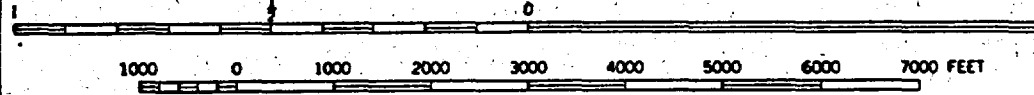
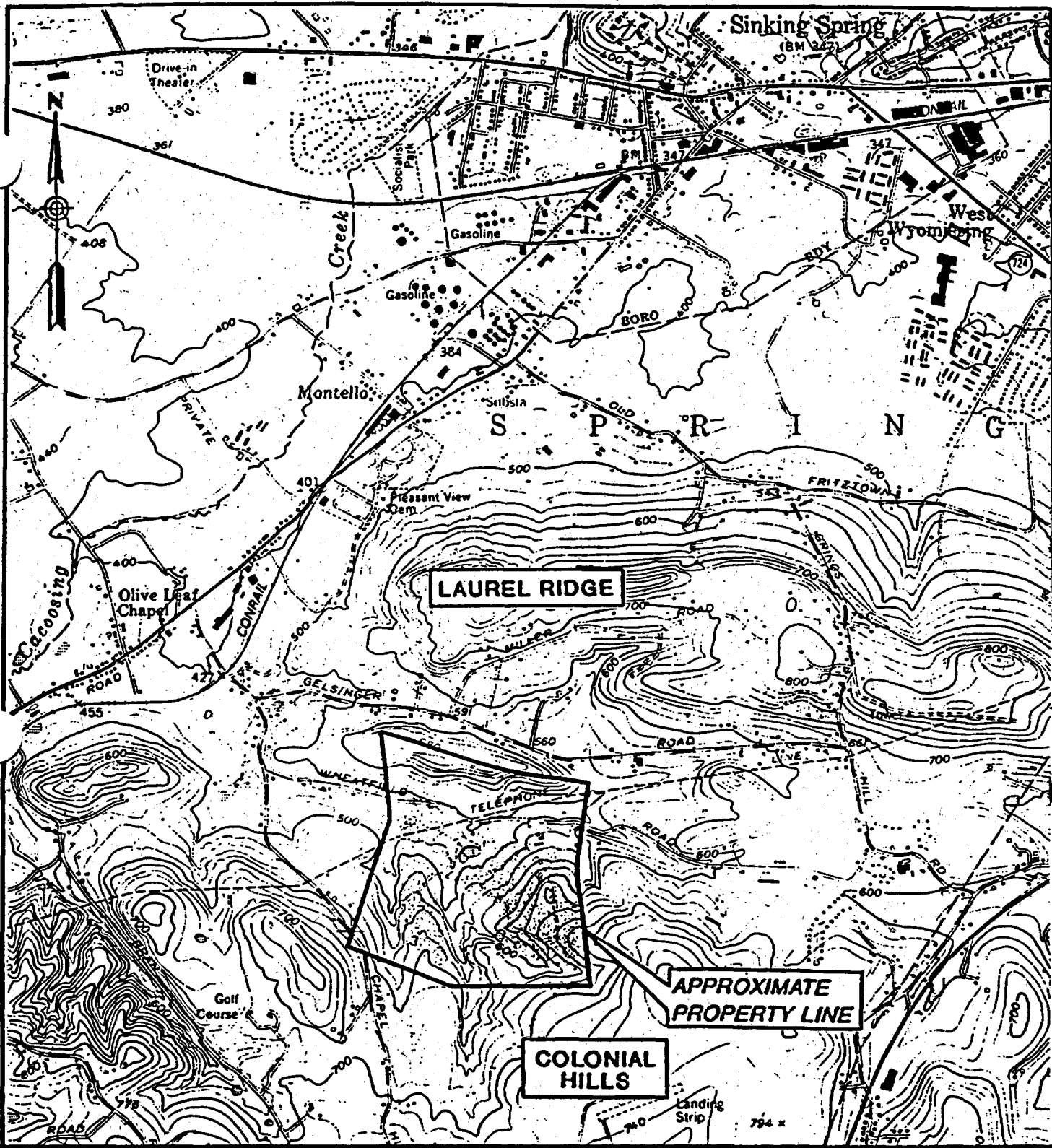

Randolph S. White, P.E.
Senior Project Manager

SAW/RSW:lrl

Attachment

z:\6773\1027\tr.wp5

AR302739



REFERENCE:
 U.S.G.S. TOPOGRAPHIC MAP, SINKING SPRING QUADRANGLE, 7.5 SERIES, PHOTOREVISED 1987.

FIG. No.:	913-6389	SCALE:	AS SHOWN
DRAWN:	EAM	DATE:	05/10/91
CHECKED:		DWG. No.:	PA21-075

OCT 27 1993

SITE LOCATION MAP

Golder Associates | **BERKS LANDFILL** | FIGURE

AR302740

158090



Commonwealth of Pennsylvania
Pennsylvania Historical and Museum Commission
Bureau for Historic Preservation
Post Office Box 1026
Harrisburg, Pennsylvania 17108-1026

JAN 27 1994

January 12, 1994

GOLDER

Stephen A. Wheeler
Golder Associates, Inc.
305 Fellowship Road, Ste. 200
Mt. Laurel, NJ 08054

TO EXPEDITE REVIEW USE
BHP REFERENCE NUMBER

Re: File No. ER 94-0309-011-A
LAND, Berks Landfill
Spring Township
Berks County

Dear Mr. Wheeler:

The Bureau for Historic Preservation has reviewed this State funded, assisted or licensed project under the authority of the Environmental Rights amendment, Article 1, Section 27 of the Pennsylvania Constitution and the Pennsylvania History Code, 37 Pa. Cons. Stat. Section 507 et seq. (1988). This review includes comments on the project's potential effect on both historic and archaeological resources.

Significant archaeological sites are located in or near your project area and others are likely to exist. These resources could be adversely affected by project activities. A Phase I archaeological survey to verify the extent of known sites and to locate other sites is needed to determine their eligibility for listing in the National Register of Historic Places. Guidelines and instructions for this phase are enclosed. The recorded sites in the area are listed below.

P.A.S.S. # 36 BK 179

In our opinion this project will have no effect on historic structures.

If you need further information in this matter please consult Andrew Wyatt at (717) 772-0923.

Sincerely,

Kurt W. Carr, Chief
Division of Archaeology &
Protection

Enclosure
KC/dks

AR302741

Golder Associates Inc.

305 Fellowship Road, Suite 200
Mt. Laurel, NJ USA 08054
Tel: (609) 273-1110
Fax (609) 273-0778



October 27, 1993

Project No.: 913-6773

- National Park Service
Mid-Atlantic Regional Office
143 South Third Street
Philadelphia, PA 19106

Attn: Ms. Patricia Bentley

RE: NATIONAL HISTORIC SITES AND NATIONAL
NATURAL LANDMARKS SEARCH

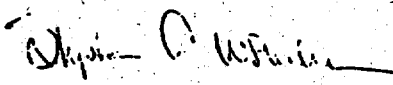
Dear Ms. Bentley:

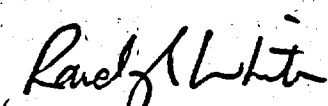
Golder Associates Inc. (Golder) requests information regarding any National Historic Sites or National Natural Landmark Sites within a five-mile radius of the Berks Landfill Site (Site). The Site is located along Wheatfield Road, Spring Township, Berks County, Pennsylvania, as shown on the enclosed Sinking Spring 7.5 minute USGS topographic quadrangle map.

If you should require additional information, please do not hesitate to contact us.

Very truly yours,

GOLDER ASSOCIATES INC.

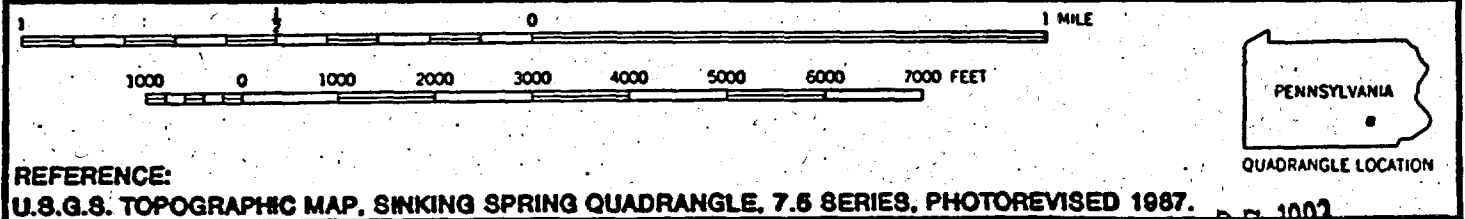
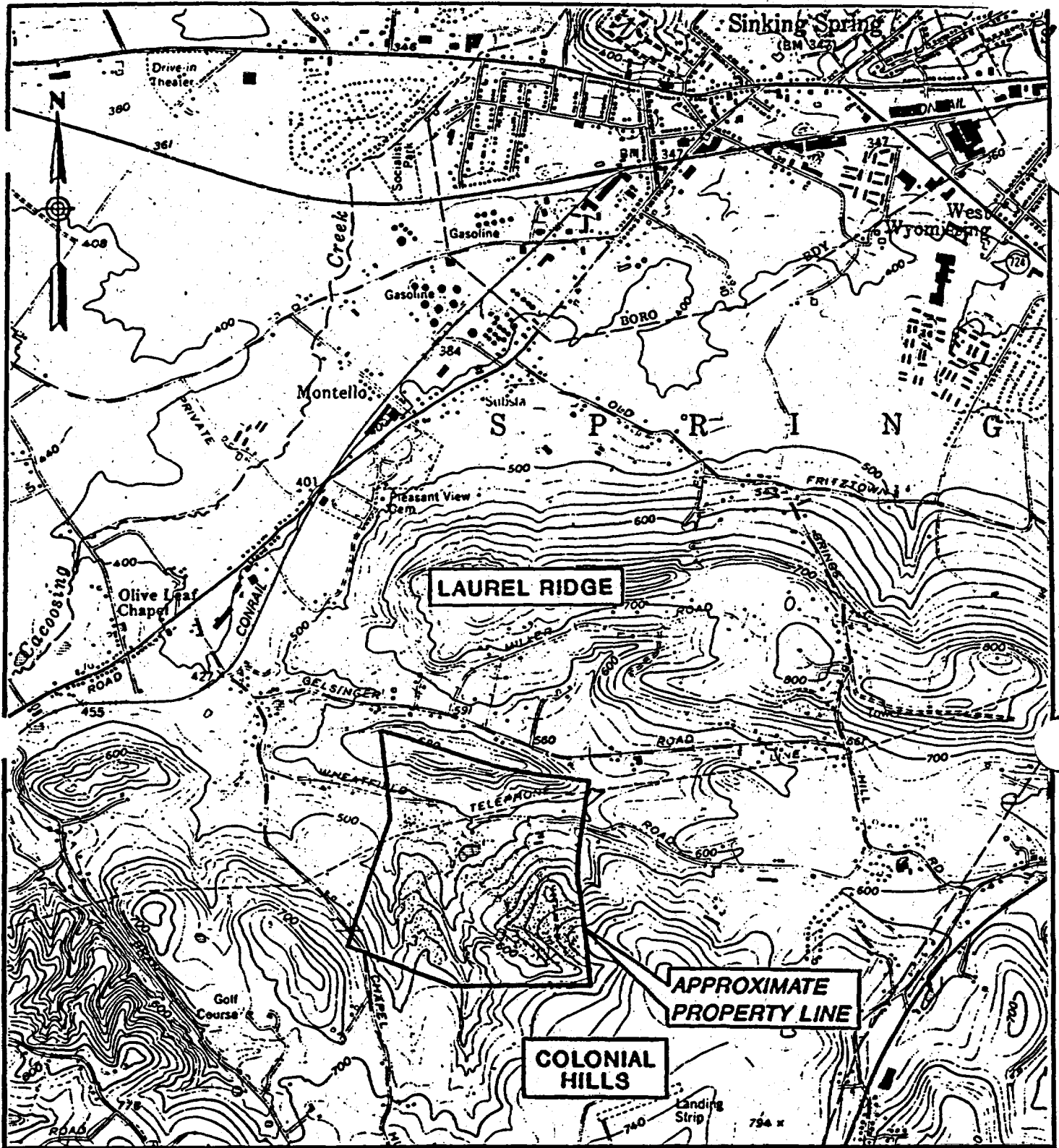

Stephen A. Wheeler
Project Environmental Scientist


Randolph S. White, P.E.
Senior Project Manager

SAW/RSW:lrl

Enclosure

z:\6773\1027\lr5.wp5



REFERENCE:
 U.S.G.S. TOPOGRAPHIC MAP, SINKING SPRING QUADRANGLE, 7.5 SERIES, PHOTOREVISED 1987.

JOB No.	913-6389	SCALE	AS SHOWN	SITE LOCATION MAP OCT 27 1993 FIGURE AR302143
DRAWN	EAM	DATE	05/10/91	
CHECKED		DWG. No.	PA21-075	
Golder Associates			BERKS LANDFILL	

10000



IN REPLY REFER TO:

United States Department of the Interior

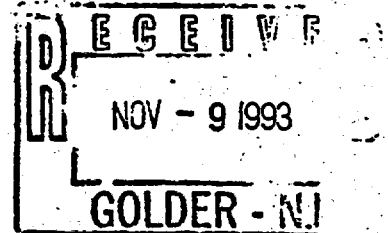


NATIONAL PARK SERVICE

Mid-Atlantic Region
143 South Third Street
Philadelphia, PA 19106

November 4, 1993

N44 (MAR-MR)



Stephen A. Wheeler
Project Environmental Scientist
and
Randolph S. White, P.E.
Senior Project Manager
Golder Associates, Inc.
305 Fellowship Road, Suite 200
Mt. Laurel, NJ 08054

Re: Project No.: 913-6773

Gentlemen:

According to the map you provided, and the area described, there are no National Natural Landmarks within a five mile radius of the project area. The closest National Natural Landmark is the Hawk Mountain Sanctuary which is approximately 30 miles north of Reading, PA.

I am forwarding a copy of your letter and enclosure to our National Historic Sites Coordinator, Bill Bolger, for his input. If I can be of further assistance, please do not hesitate to contact me. I can be reached at (215) 597-7159.

Sincerely,

Ginny Paci
National Natural Landmarks Coordinator

cc:
Bill Bolger, NHS Coord., CRM w/cy of inc.

AR302744

Golder Associates Inc.

305 Fellowship Road, Suite 200
Mt. Laurel, NJ USA 08054
Tel: (609) 273-1110
Fax (609) 273-0778



October 27, 1993

Project No.: 913-6773

U.S. Department of Interior
U.S. Fish and Wildlife Service
Suite 322
315 South Allen Street
State College, PA 16801

Attn: Mr. Charles Kulp (Field Supervisor)

RE: BERKS LANDFILL, BERKS COUNTY, PENNSYLVANIA
INFORMATION REQUEST FOR PROTECTED SPECIES AND
CULTURAL RESOURCES

Gentlemen:

Golder Associates Inc. (Golder) requests any information available regarding protected species and cultural resources within a one-mile radius of the Berks Landfill Site (Site). The Site is located along Wheatfield Road, Spring Township, Berks County, Pennsylvania, as shown on the enclosed Sinking Spring 7.5 minute USGS topographic quadrangle map.

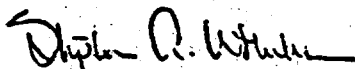
The information received will be used to address the National Environmental Protection Act (NEPA) issues in relation to the site Remedial Investigation (RI) as required under CERCLA.

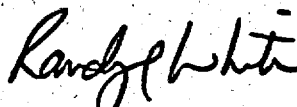
Golder has also submitted a formal request to the Pennsylvania Natural Diversity Inventory (PNDI) regarding any protected species identified or observed within a one-mile radius of the Site.

If you should require additional information, please do not hesitate to contact us.

Very truly yours,

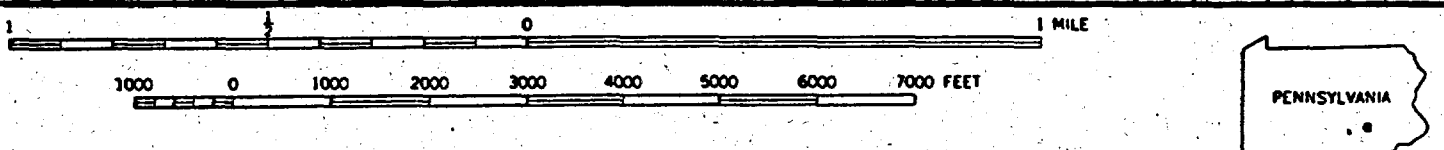
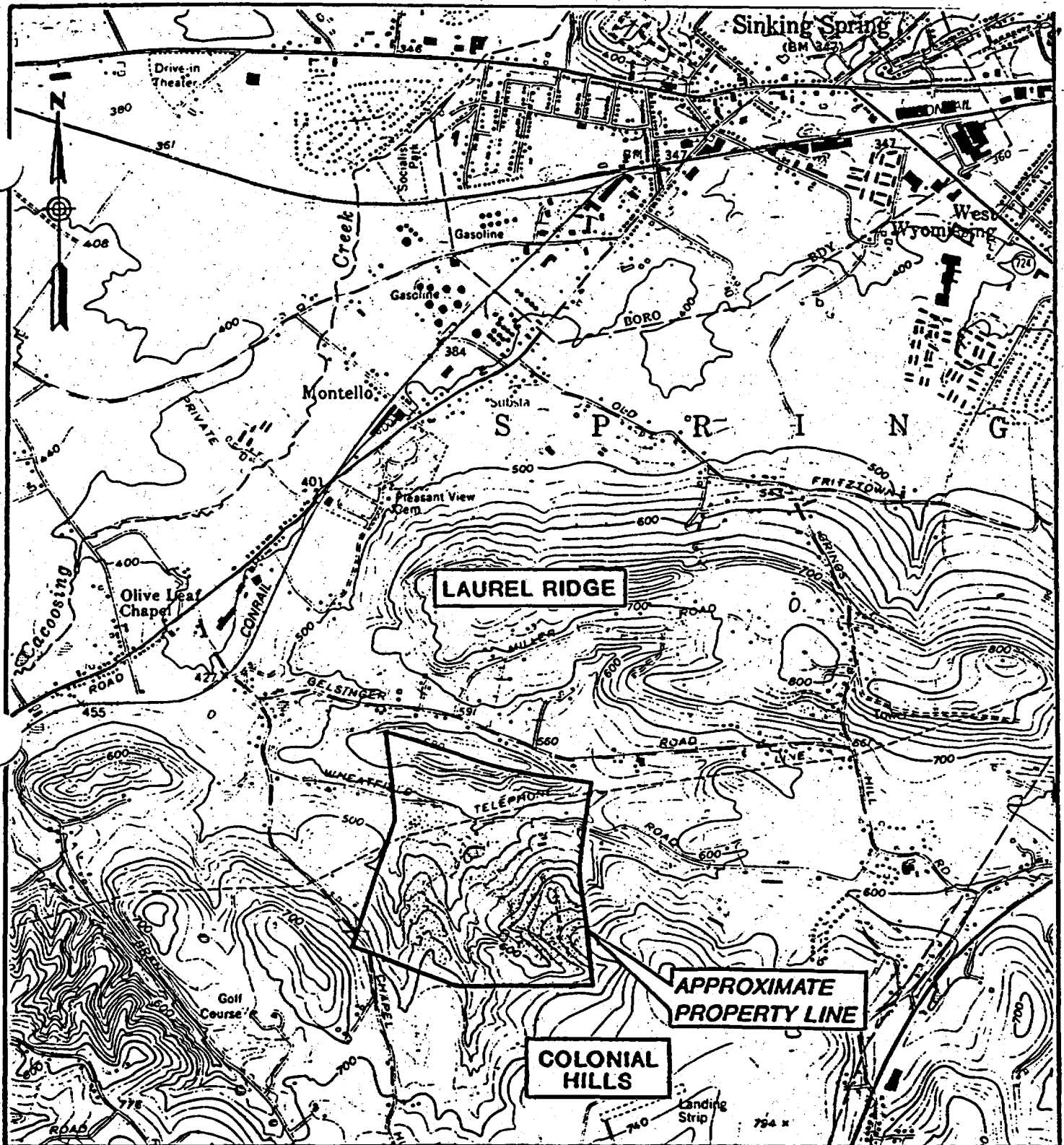
GOLDER ASSOCIATES INC.


Stephen A. Wheeler
Project Environmental Scientist


Randolph S. White, P.E.
Senior Project Manager

SAW/RSW:lrl

z:\6773\1027\tr2.wp5



REFERENCE:
 U.S.G.S. TOPOGRAPHIC MAP, SINKING SPRING QUADRANGLE, 7.6 SERIES, PHOTOREVISED 1987.

DB No.:	913-6389	SCALE:	AS SHOWN
DRAWN:	EAM	DATE:	05/10/91
CHECKED:		DWG. No.:	PA21-075

05/27/1992

SITE LOCATION MAP

Golder Associates

BERKS LANDFILL FIGURE

AR302/46

180000



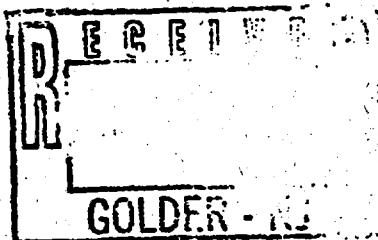
United States Department of the Interior



FISH AND WILDLIFE SERVICE
Suite 322
315 South Allen Street
State College, Pennsylvania 16801

October 29, 1993

Mr. Randolph S. White, P.E.
Senior Project Manager
Golder Associates, Inc.
305 Fellowship Road, Suite 200
Mt. Laurel, NJ 08054



Dear Mr. White:

This responds to your letter of October 27, 1993 requesting information about questions on Form D for the proposed Berks landfill site located in Spring Township, Berks County, Pennsylvania.

In response to Question 6 (Section 2, Part 1), there are no national wildlife refuges, fish hatcheries, or environmental centers operated by the Service with one mile of the site.

In response to Question 10 (Section 2, Part 1), except for occasional transient species, no federally listed or proposed threatened or endangered species under our jurisdiction are known to exist in the project impact area. Therefore, no Biological Assessment or further Section 7 consultation under the Endangered Species Act (87 Stat. 884, as amended; 16 U.S.C. 1531 et seq.) is required with the Fish and Wildlife Service. Should project plans change, or if additional information on listed or proposed species becomes available, this determination may be reconsidered. A compilation of federally listed endangered and threatened species in Pennsylvania is enclosed for your information.

Federal Candidate and State-listed Species

Candidate species are species under consideration by the Service for possible inclusion on the Federal List of Endangered and Threatened Wildlife and Plants. Because many of these species are known to have suffered population declines, the Service encourages federal agencies and other planners to consider candidate species when planning and implementing their projects.

The Pennsylvania Natural Diversity Inventory (PNDI) is maintained by the Pennsylvania Department of Environmental Resources, The Nature Conservancy and the Western Pennsylvania Conservancy. The Pennsylvania Fish and Wildlife Database is maintained by the Pennsylvania Game Commission. These databases contain the most up-to-date information about candidate and State-listed species in Pennsylvania. Requests for a PNDI review for the presence of candidate and State-listed species, as well as other natural resources of special concern, should be directed to:

Pennsylvania Department of Environmental Resources
Bureau of Forestry
Division of Forest Advisory Services
400 Market Street (MSSOB), 3rd Floor
P.O. Box 8552
Harrisburg, PA 17105-8552

AR302747

Requests for a review of the Pennsylvania Fish and Wildlife Database should be directed to:

Pennsylvania Game Commission
Bureau of Land Management
Division of Wildlife Data Base
2001 Elmerton Avenue
Harrisburg, PA 17110-9797

Should the data search reveal the presence of any candidate species on the site, the Service should be contacted to ensure that these species are not adversely affected by project activities.

Requests for information regarding State-listed endangered or threatened species should be directed to the Pennsylvania Game Commission (birds and mammals), the Pennsylvania Fish and Boat Commission (fish, reptiles, and amphibians), and the Pennsylvania Department of Environmental Resources (plants).

Section One (Exclusionary Area Criteria) questions whether the project, including any incidental earthmoving or construction activities, is located in or within 300 feet of a wetland. According to the Service's National Wetlands Inventory map for the Sinking Springs quadrangle, wetlands occur within the boundaries of the proposed project. Certain work in wetlands in Pennsylvania requires permits from the Pennsylvania Department of Environmental Resources (DER) and/or the Army Corps of Engineers (Corps). We suggest you contact the DER and the Corps at the addresses listed below for information on permit requirements.

Pennsylvania Department of
Environmental Resources
Division of Rivers and
Wetlands Conservation
P.O. Box 1467
Harrisburg, PA 17120

District Engineer, Philadelphia District
U.S. Army Corps of Engineers
100 Penn Square East
Philadelphia, PA 19107-3390

By copy of this letter, we are informing these agencies of the proposed project. You should be aware that the Service generally recommends denial of permit applications for non-water dependent fills in wetlands.

This response relates only to endangered or threatened species under our jurisdiction and a preliminary review of wetlands, based on an office review of the proposed project's location. No field inspection of the project area has been conducted by this office. Consequently, this letter is not to be construed as addressing other Service concerns under the Fish and Wildlife Coordination Act or other legislation.

If we can be of further assistance, please contact Philip Edmunds of this office at 814-234-4090.

Sincerely,



Charles J. Kulp
Supervisor

Enclosure

AR302748

FEDERALLY LISTED ENDANGERED SPECIES IN PENNSYLVANIA

<u>COMMON NAME</u>	<u>SCIENTIFIC NAME</u>	<u>STATUS*</u>	<u>DISTRIBUTION</u>
<u>FISHES</u>			
Shortnose sturgeon**	<i>Acipenser brevirostrum</i>	E	Delaware River and other Atlantic coastal waters
<u>REPTILES & AMPHIBIANS</u>			
None			
<u>BIRDS</u>			
Bald eagle	<i>Haliaeetus leucocephalus</i>	E	Entire state-recent nesting in Butler, Crawford, Dauphin, Lancaster, Pike, Tioga, York Counties
Peregrine falcon (American)	<i>Falco peregrinus anatum</i>	E	Entire state-reestablishment to former breeding range in progress
Peregrine falcon (Arctic)	<i>Falco peregrinus tundrius</i>	E	Entire state-migratory
Piping plover	<i>Charadrius melodus</i>	E	Presque Isle-no current nesting
<u>MAMMALS</u>			
Indiana bat	<i>Myotis sodalis</i>	E	Entire state
<u>MOLLUSKS</u>			
Clubshell mussel	<i>Pleurobema clava</i>	E	French Creek and Allegheny River watersheds; Clarion, Crawford, Erie, Forest, Mercer and Venango Counties
Dwarf wedge mussel	<i>Alasmidonta heterodon</i>	E	Delaware River drainage-possibly extirpated
Fanshell	<i>Cyprogenia stegaria</i>	E	Ohio River drainage-possibly extirpated
Northern riffleshell	<i>Epioblasma torulosa rangiana</i>	E	French Creek and Allegheny River watersheds; Crawford, Erie, Forest, Venango and Warren Counties
Orange pimpleback	<i>Plethobasus striatus</i>	E	Ohio River drainage-possibly extirpated
Pink mucket pearly mussel	<i>Lampsilis abrupta</i>	E	Ohio River drainage-possibly extirpated
Rough pigtoe	<i>Pleurobema plenum</i>	E	Ohio River drainage-possibly extirpated
<u>PLANTS</u>			
Northeastern bulrush	<i>Scirpus ancistrochaetus</i>	E	Current - Blair, Centre, Clinton, Cumberland, Dauphin, Franklin, Huntingdon, Lackawanna, Monroe, Union Counties. Historic - Lehigh, Northampton Counties.
Small whorled pogonia	<i>Isotria medeoloides</i>	E	Current - Centre and Venango Counties. Historic - Berks, Chester, Greene, Monroe, Montgomery, Philadelphia Counties.

* E = Endangered

** Shortnose sturgeon is under the jurisdiction of the National Marine Fisheries Service

The following is a partial list of species that no longer occur in Pennsylvania: gray wolf, eastern cougar, moose, bison, lynx, wolverine, passenger pigeon, Bachman's sparrow, common tern, lark sparrow, tiger salamander, mud sunfish, longjaw cisco, lake whitefish, butterfly mussel, American burying beetle, precious underwing moth, Karner blue butterfly, American barberry, small white lady's-slipper, eastern prairie fringed orchid, Virginia spiraea, etc, etc.

AR302749

Golder Associates Inc.

305 Fellowship Road, Suite 200
Mt. Laurel, NJ USA 08054
Tel: (609) 273-1110
Fax: (609) 273-0778



October 27, 1993

Project No.: 913-6773

National Park Service
Mid-Atlantic Regional Office
U.S. Customs House
Second & Chestnut Streets
Philadelphia, PA 19106

Attn: Ms. Katherine Stevenson

RE: NATIONAL WILD & SCENIC RIVERS AND OTHER RESOURCES SEARCH

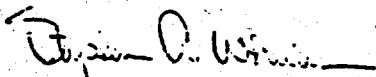
Dear Ms. Stevenson:

Golder Associates Inc. (Golder) requests information regarding National Wild and Scenic Rivers or other national resources within a five-mile radius of the Berks Landfill Site (Site). The Site is located along Wheatfield Road, Spring Township, Berks County, Pennsylvania, as shown on the enclosed Sinking Spring 7.5 minute USGS topographic quadrangle map.

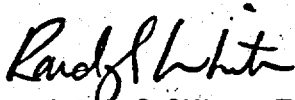
If you should require additional information, please do not hesitate to contact us.

Very truly yours,

GOLDER ASSOCIATES INC.



Stephen A. Wheeler
Project Environmental Scientist

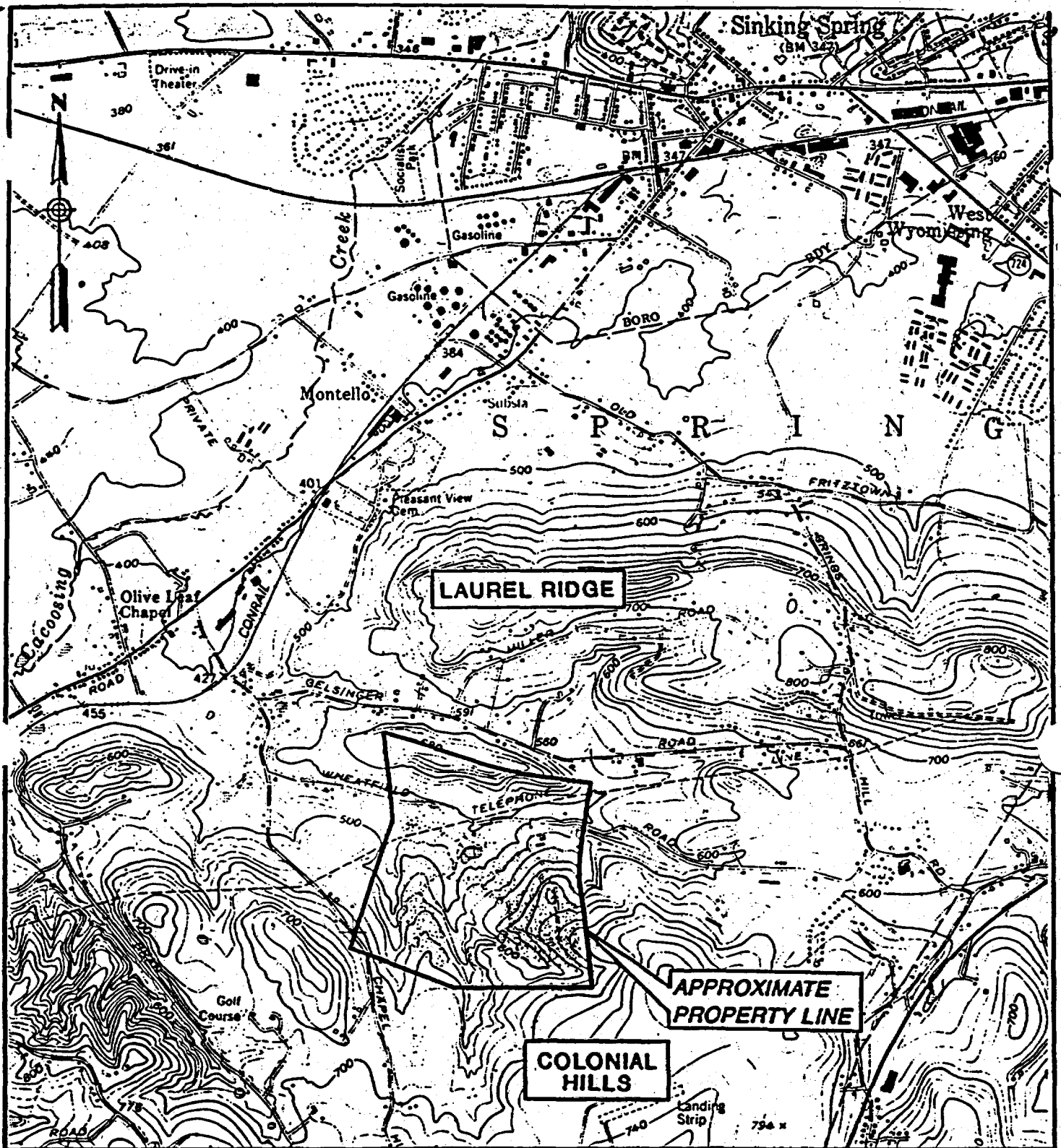


Randolph S. White, P.E.
Senior Project Manager

SAW/RSW:lrl

Enclosure

z:\6773\1027\tr4.wp5



1 0 1 MILE
 1000 0 1000 2000 3000 4000 5000 6000 7000 FEET
 PENNSYLVANIA
 QUADRANGLE LOCATION
 REFERENCE:
 U.S.G.S. TOPOGRAPHIC MAP, SINKING SPRING QUADRANGLE, 7.5 SERIES, PHOTOREVISED 1987.

JOB No.	913-6389	SCALE	AS SHOWN	SITE LOCATION MAP
DRAWN	EAM	DATE	05/10/91	
CHECKED		DWG. No.	PA21-075	
Golder Associates			BERKS LANDFILL	FIGURE

AR302751

144888



Commonwealth of Pennsylvania

R. Donald Dreese

Division of Rivers & Wetlands Conservation
Bureau of Water Resources Management
Department of Environmental Resources
3600 Vartan Way
Harrisburg, PA 17110-9333

Address:
Box 8761
Harrisburg, PA 17105-8761

(717) 541-7803

PENNSYLVANIA SCENIC RIVERS SYSTEM

System contains the following components with

Printed on recycled card stock

Accumulated Mileage

- 1: Schuylkill River - 88.4 miles - Act No. 1978-33 124.8
November 26, 1978.
Schuylkill River Extension - 31.2 miles - Act No. Pastoral, Recreat-
1988-17, March 2, 1988. ional, Modified
DER Secretary Proclamation - 5.2 miles, July 1, 1988 Recreational
Schuylkill River Greenway Association
960 Old Mill Road
Wyomissing, PA 19610
(215) 372-3916

2. Stony Creek - 16 miles - Act No. 1980-18 140.8
March 24, 1980
Pennsylvania Game Commission Wild
2001 Elmerton Ave.
Harrisburg, PA 17110-9797
(717) 787-4250

3. Lehigh River - 63.95 miles - Act No. 1982-71 204.8
April 5, 1982
Department of Environmental Resources Wild, Scenic
Bureau of State Parks
R.D. 1, Box 284
Drums, PA 18222
(717) 474-5382

4. French Creek - 42.75 miles - Act No. 1982-97 247.6
April 29, 1982
Federation of Northern Chester County Communities Scenic
R. D. 2
Pottstown, PA 19464
(215) 326-7835

5. Lick Run - 22.95 miles - Act No. 1982-324 270.6
December 17, 1982
Pennsylvania Game Commission Wild, Scenic
2001 Elmerton Ave.
Harrisburg, PA 17110-9797
(717) 787-4250

Department of Environmental Resources
Bureau of Forestry
P. O. Box 1647
Harrisburg, PA 17120
(717) 787-2703

AR302752

- | | |
|--|--|
| <p>6. Octoraro Creek - 36.5 miles - Act No. 1983-43
 October 21, 1983
 Octoraro Watershed Association
 P. O. Box 98
 Kirkwood, PA 17536
 (717) 529-2244</p> | <p>307.1
 Scenic, Pastoral</p> |
| <p>7. LeTort Spring Run - 7.6 miles - Act No. 1988-42
 March 30, 1988
 LeTort Regional Authority
 Community Center
 415 Franklin Street
 Carlisle, PA 17013
 (717) 737-5784, Ext. 25</p> | <p>314.7
 Pastoral, Modified
 Recreational</p> |
| <p>8. Bear Run - 4.3 miles - Act No. 1988-161
 December 19, 1988
 Western Pennsylvania Conservancy
 316 Fourth Avenue
 Pittsburgh, PA 15222
 (412) 288-2777</p> | <p>319.0
 Scenic</p> |
| <p>9. Tucquan Creek - 8.1 miles - Act No. 1988-161
 December 19, 1988
 Martic Hills Watershed Association
 130 Creamery Road
 Pequea, PA 17565
 (717) 284-4830</p> | <p>327.1
 Wild, Scenic</p> |
| <p>10. Lower Brandywine - 65.85 miles - Act No. 1989-7
 June 16, 1989</p> | <p>392.9
 Scenic, Pastoral</p> |

AR302753

THE NATIONAL WILD & SCENIC RIVERS SYSTEM IN PENNSYLVANIA

The National Wild & Scenic Rivers System contains the following Pennsylvania components with the managing agency:

	<u>Accumulated Mileage</u>
1. Middle Delaware River - 37.0 miles - Milford to Delaware Water Gap - September 1, 1965. National Park Service Delaware Water Gap National Recreation Area Bushkill, PA 18324 (717) 588-2435	37.0
2. Upper Delaware River - 73.4 miles - Hancock, NY to Sparrowbush, NY - November 10, 1978 National Park Service P. O. Box C Narrowsburg, PA 12764 (717) 559-7527	110.4 Scenic & Rec.

AR302754

Golder Associates Inc.

305 Fellowship Road, Suite 200
Mt. Laurel, NJ USA 08054
Tel: (609) 273-1110
Fax (609) 273-0778



November 5, 1993

Project No.: 913-6773

Pennsylvania Game Commission
Bureau of Land Management
Division of Wildlife Data Base
2001 Elmerton Avenue
Harrisburg, PA 17110-9797

RE: PENNSYLVANIA FISH AND WILDLIFE DATABASE SEARCH REQUEST
FOR THE BERKS LANDFILL SITE, BERKS COUNTY, PENNSYLVANIA

Gentlemen:

Golder Associates Inc. (Golder) requests a Pennsylvania Fish and Wildlife Database search for the Berks Landfill Site (Site), Spring Township, Berks County, Pennsylvania. Enclosed is a Site Location Map showing the Site location on the Sinking Spring, Pennsylvania, 7.5 minute USGS topographic quadrangle map.

Golder requests any information available regarding candidate and State-listed species within a one-mile radius of the Site.

If you should require additional information, please do not hesitate to contact us.

Very truly yours,

GOLDER ASSOCIATES INC.

A handwritten signature in black ink, appearing to read "Stephen A. Wheeler".

Stephen A. Wheeler
Project Environmental Scientist

SAW:mtd

z:\projects\913-6773\search

AR302755



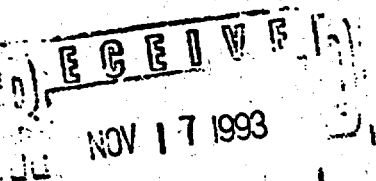
COMMONWEALTH OF PENNSYLVANIA

PENNSYLVANIA GAME COMMISSION

2001 ELMERTON AVENUE
HARRISBURG, PA 17110-9797

ADMINISTRATIVE BUREAUS:	
ADMINISTRATION	717-787-5670
AUTOMOTIVE AND PROCUREMENT DIVISION	717-787-6594
LICENSE DIVISION	717-787-2084
PERSONNEL DIVISION	717-787-7836
WILDLIFE MANAGEMENT	717-787-5529
INFORMATION & EDUCATION	717-787-6286
LAW ENFORCEMENT	717-787-5740
LAND MANAGEMENT	717-787-6818
REAL ESTATE DIVISION	717-787-6568
MANAGEMENT INFORMATION SYSTEMS	717-787-4076

November 15, 1993



Mr. Stephen A. Wheeler
Golder Associates
305 Fellowship Road
Suite 200
Mt. Laurel NJ 08054

Dear Mr. Wheeler:

In response to your request for information services, we are providing the enclosed printouts from the Pennsylvania Fish and Wildlife Data Base. This information was provided for species occurring at or near the Berks Landfill Site, Berks County, Pennsylvania.

We have no record of any endangered or threatened bird or mammal occurring at or near your project area. Additional comments concerning this data search are included on the following page.

The bill for this service is as follows:

Staff Time	9.00
Printing	1.20
<u>Mailing Cost</u>	<u>.98</u>
TOTAL	\$11.18

Please make reimbursement to the Pennsylvania Game Commission, Division of Wildlife Data Base, 2001 Elmerton Avenue, Harrisburg, PA 17110-9797.

If you have any questions or require assistance interpreting this printout, please contact Ms. Bullock at (717) 787-1570.

Very truly yours,

G.J. Grabowicz, Director
Bureau of Land Management

Pennsylvania Fish and Wildlife Data Base

The following species information was generated from the Pennsylvania Fish and Wildlife Data Base for your use in determining species likely to occur in your project area.

This information was provided upon request and should not be viewed as an official review or opinion of the Pennsylvania Game Commission. Species lists generated for this request should be viewed as likely or probable occurrence lists that might warrant further investigation. These lists are based on known, documented species occurrence within the counties, watershed, land use, and/or habitat types specified in your request.

Information pertaining to aquatic vertebrates and invertebrates contained in these lists is based solely on literature sources and expert opinion. Use of the aquatic species information contained in this report should be coordinated with the Pennsylvania Fish and Boat Commission for compliance with their standards and data sources.

This report does not contain information on plants. For plant species information in your project area, please contact the Bureau of Forestry, Pennsylvania Department of Environmental Resources.

The Pennsylvania Game Commission considers wetlands critical and unique wildlife habitat. If your proposed project is in the vicinity of wetlands, streams, rivers, lakes, or other bodies of water, please be aware that any impact to these areas requires a permit from the U.S. Army Corps of Engineers and the Bureau of Dams and Waterway Management, Pennsylvania Department of Environmental Resources.

Note: Bird species occurrence is based upon recorded sightings and may not imply nesting activity or year-round residence.

Contents

- A. Project Area Endangered and Threatened Species List
- B. Project Area Potential Special Concern Species List
- C. Land Use/Cover Type Table for Potential Special Concern Species
- D. MASTER LIST - Project Area Species List

AR302758

Pennsylvania Fish and Wildlife Data Base
LIST A: Endangered and Threatened Species
** Berks Landfill Site **
Sinking Spring Quadrangle
15 NOV 1993

Note: The purpose of the following list is to identify endangered or threatened species which occur or are likely to occur on a designated site. We have record of the following species occurring in or near your project area. Their occurrence may depend on season, habitat type, and individual movements or migration patterns. Field surveys may be required to determine whether these species exist on your project area.

Species.ID Common Name..... Scientific Name.....

ZERO Records Listed

Pennsylvania Fish and Wildlife Data Base
LIST B: Potential Special Concern Species
(Includes Accidental and Migrant Species)

** Berks Landfill Site **

Berks County

15 NOV 1993

Note: The purpose of the following list is to identify endangered, threatened, and special concern species which may potentially occur within a designated area. This list includes species which may exist on your project area as well as migrating and accidental species. This information is based on records of these animals inhabiting specific habitat types within Berks County.

Status.....	No. of Species Listed
PA / Fed Endangered	2
PA Endangered	6
Fed Endangered	1
PA Threatened	3
Candidate Species	32
Species Listed:	44

AR302760

Pennsylvania Fish and Wildlife Data Base
 LIST B: Potential Special Concern Species
 (Includes Accidental and Migrant Species)
 ** Berks Landfill Site **
 Berks County
 15 NOV 1993

Common Name.....	Scientific Name.....	Status.....
Eagle, Bald	<i>Haliaeetus leucocephalus</i>	PA / Fed Endangered
Falcon, Peregrine	<i>Falco peregrinus</i>	PA / Fed Endangered
Turtle, Bog	<i>Clemmys muhlenbergi</i>	PA Endangered
Osprey	<i>Pandion haliaetus</i>	PA Endangered
Owl, Short-eared	<i>Asio flammeus</i>	PA Endangered
Rail, King	<i>Rallus elegans</i>	PA Endangered
Tern, Black	<i>Chlidonias niger</i>	PA Endangered
Shrew, Least	<i>Cryptotis parva</i>	PA Endangered
Tern, Least	<i>Sterna antillarum</i>	Fed Endangered
Bittern, American	<i>Botaurus lentiginosus</i>	PA Threatened
Flycatcher, Yellow-bellied	<i>Empidonax flaviventris</i>	PA Threatened
Sandpiper, Upland	<i>Bartramia longicauda</i>	PA Threatened
Harrier, Northern	<i>Circus cyaneus</i>	Candidate - At Risk
Owl, Common Barn	<i>Tyto alba</i>	Candidate - At Risk
Snipe, Common	<i>Gallinago gallinago</i>	Candidate - At Risk
Sparrow, Henslow's	<i>Ammodramus henslowii</i>	Candidate - At Risk
Warbler, Prothonotary	<i>Protonotaria citrea</i>	Candidate - At Risk
Bobcat	<i>Felis rufus</i>	Candidate - At Risk
Hare, Snowshoe	<i>Lepus americanus</i>	Candidate - At Risk
Coot, American	<i>Fulica americana</i>	Candidate - Rare
Goshawk, Northern	<i>Accipiter gentilis</i>	Candidate - Rare
Grebe, Pied-billed	<i>Podilymbus podiceps</i>	Candidate - Rare
Grosbeak, Blue	<i>Guiraca caerulea</i>	Candidate - Rare
Tanager, Summer	<i>Piranga rubra</i>	Candidate - Rare
Teal, Green-winged	<i>Anas crecca</i>	Candidate - Rare
Thrush, Swainson's	<i>Catharus ustulatus</i>	Candidate - Rare
Wren, Marsh	<i>Cistothorus palustris</i>	Candidate - Rare
Bat, Silver-haired	<i>Lasionycteris noctivagans</i>	Candidate - Rare
Bobwhite, Northern	<i>Colinus virginianus</i>	Candidate - Undeterm
Crossbill, Red	<i>Loxia curvirostra</i>	Candidate - Undeterm

Pennsylvania Fish and Wildlife Data Base
 LIST B: Potential Special Concern Species
 (Includes Accidental and Migrant Species)
 ** Berks Landfill Site **
 Berks County
 15 NOV 1993

Common Name.....	Scientific Name.....	Status.....
Dickcissel	<i>Spiza americana</i>	Candidate - Undeterm
Duck, Ruddy	<i>Oxyura jamaicensis</i>	Candidate - Undeterm
Egret, Cattle	<i>Bubulcus ibis ibis</i>	Candidate - Undeterm
Gadwall	<i>Anas strepera</i>	Candidate - Undeterm
Nighthawk, Common	<i>Chordeiles minor</i>	Candidate - Undeterm
Owl, Long-eared	<i>Otus asio</i>	Candidate - Undeterm
Owl, Northern Saw-whet	<i>Aegolius acadicus</i>	Candidate - Undeterm
Pintail, Northern	<i>Anas acuta</i>	Candidate - Undeterm
Shoveler, Northern	<i>Anas clypeata</i>	Candidate - Undeterm
Whip-poor-will	<i>Caprimulgus vociferus</i>	Candidate - Undeterm
Wigeon, American	<i>Anas americana</i>	Candidate - Undeterm
Wren, Seminole	<i>Lasiurus seminolus</i>	Candidate - Undeterm
Bowfin	<i>Amia calva</i>	Candidate Species
Rattlesnake, Timber	<i>Crotalus horridus</i>	Candidate Species

PRIME FARMLAND

Survey Area- BERKS COUNTY, PENNSYLVANIA

Map Symbol	Prime Farmland Code	Soil Map Unit Name
AgB	1	ALLENWOOD GRAVELLY SILT LOAM, 2 TO 8 PERCENT SLOPES
AsB2	1	ATHOL SILT LOAM, 3 TO 8 PERCENT SLOPES, MODERATELY ERODED
AsC2	S	ATHOL SILT LOAM, 8 TO 15 PERCENT SLOPES, MODERATELY ERODED
AsC3	S	ATHOL SILT LOAM, 8 TO 15 PERCENT SLOPES, SEVERELY ERODED
Au	S	ATKINS SILT LOAM
BeB2	1	BEDINGTON SHALY SILT LOAM, 3 TO 8 PERCENT SLOPES, MODERATELY ERODED
BeC2	S	BEDINGTON SHALY SILT LOAM, 8 TO 15 PERCENT SLOPES, MODERATELY ERODED
BkA2	S	BERKS SHALY SILT LOAM, 0 TO 3 PERCENT SLOPES, MODERATELY ERODED
BkB2	S	BERKS SHALY SILT LOAM, 3 TO 8 PERCENT SLOPES, MODERATELY ERODED
BkC2	S	BERKS SHALY SILT LOAM, 8 TO 15 PERCENT SLOPES, MODERATELY ERODED
BlB	1	BIRDSBORO SILT LOAM, 2 TO 10 PERCENT SLOPES
BmB	1	BIRDSBORO-DUFFIELD SILT LOAMS, 3 TO 10 PERCENT SLOPES
Bo	S	BOWMANVILLE SILT LOAM
BrB2	S	BRANDYWINE CHANNERY LOAM, 3 TO 8 PERCENT SLOPES, MODERATELY ERODED
BrC2	S	BRANDYWINE CHANNERY LOAM, 8 TO 15 PERCENT SLOPES, MODERATELY ERODED
BsB	1	BRECKNOCK CHANNERY SILT LOAM, 3 TO 8 PERCENT SLOPES
BsC2	S	BRECKNOCK CHANNERY SILT LOAM, 8 TO 15 PERCENT SLOPES, MODERATELY ERODED
BvB	1	BUCHANAN GRAVELLY LOAM, 3 TO 8 PERCENT SLOPES
ChB2	1	CHESTER CHANNERY SILT LOAM, 3 TO 8 PERCENT SLOPES, MODERATELY ERODED
ChC2	S	CHESTER CHANNERY SILT LOAM, 8 TO 15 PERCENT SLOPES, MODERATELY ERODED
ChC3	S	CHESTER CHANNERY SILT LOAM, 8 TO 15 PERCENT SLOPES, SEVERELY ERODED
CoA	1	COMLY SILT LOAM, 0 TO 3 PERCENT SLOPES
CoB2	1	COMLY SILT LOAM, 3 TO 8 PERCENT SLOPES, MODERATELY ERODED
DfA	1	DUFFIELD SILT LOAM, 0 TO 3 PERCENT SLOPES
DfB2	1	DUFFIELD SILT LOAM, 3 TO 8 PERCENT SLOPES, MODERATELY ERODED
DfC2	S	DUFFIELD SILT LOAM, 8 TO 15 PERCENT SLOPES, MODERATELY ERODED
DhC3	S	DUFFIELD AND HAGERSTOWN SOILS, 8 TO 15 PERCENT SLOPES, SEVERELY ERODED
EcB2	1	EDGEMONT CHANNERY LOAM, 3 TO 8 PERCENT SLOPES, MODERATELY ERODED
EcC2	S	EDGEMONT CHANNERY LOAM, 8 TO 15 PERCENT SLOPES, MODERATELY ERODED
FoA	1	FOGELSVILLE SILT LOAM, 0 TO 3 PERCENT SLOPES
FoB2	1	FOGELSVILLE SILT LOAM, 3 TO 8 PERCENT SLOPES, MODERATELY ERODED

AR302763

PRIME FARMLAND

Survey Area- BERKS COUNTY, PENNSYLVANIA

Map Symbol	Prime Farmland Code	Soil Map Unit Name
FoC2	S	FOGELSVILLE SILT LOAM, 8 TO 15 PERCENT SLOPES, MODERATELY ERODED
GlA	1	GLENVILLE SILT LOAM, 0 TO 3 PERCENT SLOPES
GlB2	1	GLENVILLE SILT LOAM, 3 TO 8 PERCENT SLOPES, MODERATELY ERODED
HaB2	1	HAGERSTOWN SILT LOAM, 3 TO 8 PERCENT SLOPES, MODERATELY ERODED
HaC2	S	HAGERSTOWN SILT LOAM, 8 TO 15 PERCENT SLOPES, MODERATELY ERODED
KlB2	S	KLINESVILLE SHALY SILT LOAM, 3 TO 8 PERCENT SLOPES, MODERATELY ERODED
KlC2	S	KLINESVILLE SHALY SILT LOAM, 8 TO 15 PERCENT SLOPES, MODERATELY ERODED
LaB2	1	LAIDIG CHANNERY LOAM, 3 TO 8 PERCENT SLOPES, MODERATELY ERODED
LaC2	S	LAIDIG CHANNERY LOAM, 8 TO 15 PERCENT SLOPES, MODERATELY ERODED
LhA	1	LEHIGH SILT LOAM, 0 TO 3 PERCENT SLOPES
LhB2	1	LEHIGH SILT LOAM, 3 TO 8 PERCENT SLOPES, MODERATELY ERODED
LhC3	S	LEHIGH SILT LOAM, 8 TO 15 PERCENT SLOPES, SEVERELY ERODED
LrB2	1	LEWISBERRY GRAVELLY SANDY LOAM, 3 TO 8 PERCENT SLOPES, MODERATELY ERODED
LrC2	S	LEWISBERRY GRAVELLY SANDY LOAM, 8 TO 15 PERCENT SLOPES, MODERATELY ERODED
LrC3	S	LEWISBERRY GRAVELLY SANDY LOAM, 8 TO 15 PERCENT SLOPES, SEVERELY ERODED
Lt	1	LINDSIDE SILT LOAM
LzB2	S	LITZ SHALY SILT LOAM, 3 TO 8 PERCENT SLOPES, MODERATELY ERODED
LzC2	S	LITZ SHALY SILT LOAM, 8 TO 15 PERCENT SLOPES, MODERATELY ERODED
Ml	S	MELVIN SILT LOAM
MrA3	1	MURRILL GRAVELLY CLAY LOAM, 0 TO 3 PERCENT SLOPES, SEVERELY ERODED
MuA	1	MURRILL GRAVELLY LOAM, 0 TO 3 PERCENT SLOPES
MuB2	1	MURRILL GRAVELLY LOAM, 3 TO 8 PERCENT SLOPES, MODERATELY ERODED
MuC2	S	MURRILL GRAVELLY LOAM, 8 TO 15 PERCENT SLOPES, MODERATELY ERODED
✓NaC3	S	NESHAMINY SILTY CLAY LOAM, 8 TO 15 PERCENT SLOPES, SEVERELY ERODED
NeB2	1	NESHAMINY SILT LOAM, 3 TO 8 PERCENT SLOPES, MODERATELY ERODED
NeC2	S	NESHAMINY SILT LOAM, 8 TO 15 PERCENT SLOPES, MODERATELY ERODED
PeB2	1	PENN SHALY SOILS, 3 TO 8 PERCENT SLOPES, MODERATELY ERODED
PeC2	S	PENN SHALY SOILS, 8 TO 15 PERCENT SLOPES, MODERATELY ERODED
Ph	S	PHILO LOAM, COAL OVERWASH

AR302764

PRIME FARMLAND

Survey Area- BERKS COUNTY, PENNSYLVANIA

Map Symbol	Prime Farmland Code	Soil Map Unit Name
P1	1	PHILO SILT LOAM
Po	1	POPE SILT LOAM
RaB	1	RARITAN SILT LOAM, 0 TO 5 PERCENT SLOPES
ReA	1	READINGTON SILT LOAM, 0 TO 3 PERCENT SLOPES
ReB2	S	READINGTON SILT LOAM, 3 TO 8 PERCENT SLOPES, MODERATELY ERODED
RIa2	S	REAVILLE SHALY SILT LOAM, 0 TO 3 PERCENT SLOPES, MODERATELY ERODED
RIb2	S	REAVILLE SHALY SILT LOAM, 3 TO 8 PERCENT SLOPES, MODERATELY ERODED
Ro	1	ROWLAND SILT LOAM
RyA2	1	RYDER SILT LOAM, 0 TO 3 PERCENT SLOPES, MODERATELY ERODED
RyB2	1	RYDER SILT LOAM, 3 TO 8 PERCENT SLOPES, MODERATELY ERODED
RyC3	S	RYDER SILT LOAM, 8 TO 15 PERCENT SLOPES, SEVERELY ERODED
WaA2	1	WASHINGTON SILT LOAM, 0 TO 3 PERCENT SLOPES, MODERATELY ERODED
WaB2	1	WASHINGTON SILT LOAM, 3 TO 8 PERCENT SLOPES, MODERATELY ERODED
WaC2	S	WASHINGTON SILT LOAM, 8 TO 15 PERCENT SLOPES, MODERATELY ERODED
WcA	1	WATSON SILT LOAM, 0 TO 3 PERCENT SLOPES
WeB2	S	WEIKERT-BERKS SHALY SILT LOAMS, 3 TO 8 PERCENT SLOPES, MODERATELY ERODED
WeC2	S	WEIKERT-BERKS SHALY SILT LOAMS, 8 TO 15 PERCENT SLOPES, MODERATELY ERODED
WsA	1	WILTSHIRE SILT LOAM, 0 TO 3 PERCENT SLOPES
WsB2	1	WILTSHIRE SILT LOAM, 3 TO 8 PERCENT SLOPES, MODERATELY ERODED

Prime Farmland

Code Description

- 1 All areas are Prime Farmland
- S Map Units that are Statewide Important Farmland

AR302765

APPENDIX P
LABORATORY ANALYTICAL RESULTS

AR302766

GROUNDWATER

AR302767

APPENDIX P**SUMMARY OF DATA PRESENTATION**

For each phase of sampling, (Phase 1A, 1B [including residential well sampling], and 1C [soil]), analytical data was received from the laboratory on diskette and in hard copy format. The hard copy of the data was validated, and a validation narrative was written. A database was created using the data provided by the laboratory on diskette. Laboratory qualifiers in the database were revised according to validator's qualifiers (defined in Table P-1, attached to this Appendix). All validated analytical data (sample results [including re-analyses and field duplicates] and validated qualifiers) were tabulated in a format patterned after the data summary forms in the USEPA Region III data validation guidelines. A complete copy of the data package and the validation narrative for each phase were submitted to USEPA during the course of the Remedial Investigation (RI).

This appendix of the RI Report contains the Berks Landfill RI database which includes all validated analytical data for each phase of sampling but does not include USEPA split sampling results. This appendix includes primary and re-analysis results for samples which may have required more than one analysis due to quality control criteria or the need for a dilution. The data are arranged by sampling medium (groundwater, residential well, soil, leachate, surface water, and sediment). Within each medium group, the results are arranged by 1) analytical fraction--volatile; semivolatile; pesticide and PCB; total inorganics (metals and cyanide); dissolved inorganics (metals); and wet chemistry and 2) by sampling phase.

The individual databases created for each phase were copied into one large database. From this large database, a subset of analytical results consisting of constituents considered to be present in each medium subgroup (e.g., background groundwater; see Appendix Q of the RI report, "Explanation and Notes," Table Q-A for list of subgroups) was defined and tabulated. This subset database was created to provide a clearer

presentation of constituents detected in the various media studied. As explained in Appendix Q, "Explanation and Notes," within a medium subgroup, a constituent was considered to be present if at least one detected concentration (data qualified with an A, J, K, or L) for that constituent was reported. Data for each constituent determined to be present in a subgroup has been tabulated for each sample point associated with the subgroup. The tables are arranged by analytical fraction. Tables of constituents considered to be present are included in the RI report as Appendix Q.

Calculations then were performed on the data in Appendix Q to determine frequency of detection, range of detected concentrations, and mean concentration for the constituents considered to be present. Tables summarizing the calculations are included in Section 4.0 of the RI report. These data summary tables in Section 4.0 (Tables 4-2 through 4-22) are based upon data presented in Appendix Q. The tables summarize frequency of detection, range of concentration, and mean concentration of each detected constituent in each media subgroup. The summaries were prepared in accordance with procedures detailed in Appendix R of this report. Anomalies in data and the manner in which anomalies were handled are also discussed in Appendix R.

After calculations were performed on validated analytical data for total and dissolved inorganics in background groundwater and on-site groundwater in Appendix Q of the RI report, the inorganic results were compared using the Tolerance Interval Approach ("Statistical Analysis of Groundwater Monitoring Data at RCRA Facilities, USEPA, 1989), and the Wilcoxin Two Sample Test (Sokal and Rohlf, 1969). The comparisons are included in this report as Appendix T of the RI report.

Chemistry figures displaying detected results for all sampling phases by medium and analytical group (e.g., "SHALLOW GROUNDWATER MONITORING WELL AND RESIDENTIAL WELL SAMPLE ORGANIC ANALYSES RESULTS") were produced and are included in Section 4.0 of the RI report. Due to the large number of analytical values reported for each sample point, Golder chose a "representative summary data set" for

presentation on summary figures (Figures 4-1 to 4-10) for those sample points which had re-analyses or dilutions. A conservative approach was used to select the "representative summary data set." In reviewing the sample and re-analysis data, the following steps were used in selecting this "representative summary data set":

1. If the analyte was detected in both analyses and qualified (A, J, K, or L), the higher result was chosen regardless of the qualifier.
2. If one result was deemed usable (A, J, K, or L) and the other result was deemed unusable (R) or tentatively identified (N), the usable result was chosen.
3. If the analyte was only detected in one analysis, the detected result was chosen over the non-detected result.

In situations where a field duplicate was collected, both the primary and the field duplicate sample results are presented on the figure (Figures 4-1 through 4-10). Split sampling results are not presented on the figures but have been included in Attachment 2. Exceptions to this approach are noted in Appendix R for anomalous situations.

Figures 4-1 through 4-5 present the "representative summary data set" for Phase 1A and 1B groundwater and residential sampling points and include field duplicate sampling results. Figures 4-6 and 4-7 present the "representative summary data set" for Phase 1A, 1B, and 1C soil and include field duplicate sampling results. Figure 4-8 contains the "representative summary data set" for Phase 1A and 1B leachate and includes field duplicate sampling results. Figures 4-9 and 4-10 present the "representative summary data set" for Phase 1A and 1B surface water and sediments and include field duplicate sampling results.

Discussion of chemistry data in the RI text is based upon the data presented on the figures except where specifically noted otherwise because the figures provide spatial representations of the chemicals detected in environmental media. Data for samples having field duplicates are discussed in terms of average values calculated from the primary and field duplicate results. In some instances, two results are presented in the text for a particular well because these are the results from the Phase 1A and Phase 1B sampling events.

VOLATILE ORGANIC COMPOUNDS

Phase 1A
Phase 1B

AR302771

SUMMARY OF CLP ORGANIC ANALYSES - PHASE 1A
BERKS LANDFILL REMEDIAL INVESTIGATION
BERKS COUNTY, PENNSYLVANIA

Matrix: Groundwater Well

Volatile Organics

CRQL	Parameter	Sample Point GR180/1A			Sample Point GV-16/1A			Sample Point GV-C5/1A			Sample Point GV185/1A		
		SQL	Result	Qual	SQL	Result	Qual	SQL	Result	Qual	SQL	Result	Qual
10	Acetone	10	10	U	10	32	B	10	10	U	170	U	
10	Benzene	10	10	U	10	13	A	10	10	J	170	U	
10	Bromodichloromethane	10	10	U	10	10	U	10	10	U	170	U	
10	Bromoform	10	10	U	10	10	U	10	10	U	170	U	
10	Bromomethane	10	10	U	10	10	U	10	10	U	170	U	
10	2-Butanone	10	10	U	10	39	A	10	10	U	170	U	
10	Carbon Disulfide	10	10	U	10	10	U	10	10	U	170	U	
10	Carbon Tetrachloride	10	10	U	10	10	U	10	10	U	170	U	
10	Chlorobenzene	10	10	U	10	4	J	10	10	U	170	U	
10	Chloroethane	10	10	U	10	10	U	10	10	U	170	U	
10	Chloroform	10	10	U	10	10	U	10	10	U	170	U	
10	Dibromochloromethane	10	10	U	10	10	U	10	10	U	170	U	
10	1,1-Dichloroethane	10	10	U	10	10	U	10	10	U	170	U	
10	1,2-Dichloroethane	10	10	U	10	10	U	10	10	U	170	U	
10	1,1-Dichloroethene	10	10	U	10	10	U	10	10	U	170	U	
10	Total 1,2-Dichloroethene	10	10	U	10	10	U	10	10	U	170	U	
10	1,2-Dichloropropane	10	10	U	10	10	U	10	10	U	170	U	
10	cis-1,3-Dichloropropene	10	10	U	10	10	U	10	10	U	170	U	
10	trans-1,3-Dichloropropene	10	10	U	10	10	U	10	10	U	170	U	
10	Ethylbenzene	10	10	U	10	23	A	10	10	U	170	U	
10	2-Hexanone	10	10	U	10	10	U	10	10	U	170	U	
10	4-Methyl-2-Pentanone	10	10	U	10	10	U	10	10	U	170	U	
10	Methylene Chloride	10	10	U	10	10	U	10	10	U	170	U	
10	Styrene	10	10	U	10	10	U	10	10	U	170	U	
10	1,1,2,2-Tetrachloroethane	10	10	U	10	10	U	10	10	U	170	U	
10	Tetrachloroethene	10	10	U	10	10	U	10	10	U	170	U	
10	Toluene	10	10	U	10	10	U	10	10	U	170	U	
10	1,1,1-Trichloroethane	10	10	U	10	10	U	10	10	U	170	U	
10	1,1,2-Trichloroethane	10	10	U	10	10	U	10	10	U	170	U	
10	Trichloroethene	10	10	U	10	10	U	10	10	U	170	U	
10	Vinyl Chloride	10	10	U	10	10	U	10	10	J	3100	A	
10	Total Xylenes	10	10	U	10	110	A	10	10	U	180	A	
10		10	10	U	10	10	U	10	10	U	170	U	

Notes:
All units are ug/l
Sample Quantitation Limit (SQL) is CRQL multiplied by the dilution factor and adjusted for percent moisture as appropriate.
The Qual column indicates the qualifier applied to the result following data validation (see below).

Qualifiers:
A - Acceptable (Quantitative) Data
J - Estimated (Semi-quantitative) Data
B - Present in associated blanks (Semi-quantitative Data)
K - Biased High (Semi-quantitative) Data
L - Biased Low (Semi-quantitative Data)
UJ - Not Detected/Estimated Data
U - Not Detected/Quantitative Data
UL - Not Detected/Biased Low
M - Tentative Identification
R - Unusable Data
Q - Not Applicable

AR302772

SUMMARY OF CLP ORGANICS ANALYSES - PHASE 1A
BERKS LANDFILL REMEDIAL INVESTIGATION
BERKS COUNTY, PENNSYLVANIA

Matrix: Groundwater Well

Volatile Organics

CPL	Parameter	Sample Point GU185/F0		Sample Point GU350/1A		Sample Point GU301/1A		Sample Point GU301/1A	
		SQL	Result	SQL	Result	SQL	Result	SQL	Result
10	Acetone	200	200	200	200	10	10	10	10
10	Benzene	200	200	200	200	10	10	10	10
10	Bromodichloromethane	200	200	200	200	10	10	10	10
10	Bromoform	200	200	200	200	10	10	10	10
10	Bromomethane	200	200	200	200	10	10	10	10
10	2-Butanone	200	200	200	200	10	10	10	10
10	Carbon Disulfide	200	200	200	200	10	10	10	10
10	Carbon Tetrachloride	200	200	200	200	10	10	10	10
10	Chlorobenzene	200	200	200	200	10	10	10	10
10	Chloroethane	200	200	200	200	10	10	10	10
10	Chloroform	200	200	200	200	10	10	10	10
10	Chloromethane	200	200	200	200	10	10	10	10
10	Dibromochloroethane	200	200	200	200	10	10	10	10
10	1,1-Dichloroethane	200	200	200	200	10	10	10	10
10	1,2-Dichloroethane	200	200	200	200	10	10	10	10
10	1,1-Dichloroethene	200	200	200	200	10	10	10	10
10	Total 1,2-Dichloroethene	200	2100	200	3700	10	62	10	10
10	1,2-Dichloropropane	200	200	200	200	10	10	10	10
10	cis-1,3-Dichloropropene	200	200	200	200	10	10	10	10
10	trans-1,3-Dichloropropene	200	200	200	200	10	10	10	10
10	Ethylbenzene	200	200	200	200	10	10	10	10
10	2-Hexanone	200	200	200	200	10	10	10	10
10	4-Methyl-2-Pentanone	200	200	200	200	10	10	10	10
10	Methylene Chloride	200	130	200	200	10	11	10	3
10	Styrene	200	200	200	200	10	10	10	10
10	1,1,2-Tetrachloroethane	200	200	200	200	10	10	10	10
10	Tetrachloroethene	200	200	200	200	10	10	10	10
10	Toluene	200	200	200	200	10	10	10	10
10	1,1,1-Trichloroethane	200	200	200	200	10	10	10	10
10	1,1,2-Trichloroethane	200	200	200	200	10	10	10	10
10	Trichloroethene	200	3100	200	2000	10	4	10	10
10	Vinyl Chloride	200	170	200	370	10	13	10	10
10	Total Xylenes	200	200	200	200	10	10	10	10

Notes:
 All units are ug/l
 Sample Quantitation Limit (SQL) is CROL multiplied by the dilution factor and adjusted for percent moisture as appropriate.
 The Qual column indicates the qualifier applied to the result following data validation (see below).

Qualifiers:.....
 A - Acceptable (Quantitative) Data K - Biased High (Semi-quantitative) Data UJ - Not Detected/Estimated Data M - Tentative Identification
 J - Estimated (Semi-quantitative) Data L - Biased Low (Semi-quantitative) Data U - Not Detected/Quantitative Data R - Unusable Data
 B - Present in associated blanks (Semi-quantitative Data) UL - Not Detected/Biased Low O - Not Applicable

AR302773

913
 SUMMARY OF CLP ORGANICS ANALYSES - PHASE 1A
 BERKS LANDFILL REMEDIAL INVESTIGATION
 BERKS COUNTY, PENNSYLVANIA

Matrix: Groundwater Well

Volatiles Organics

CROL	Parameter	Sample Point GMG02/1A			Sample Point GMG03/1A			Sample Point		
		SQL	Result	Qual	SQL	Result	Qual	SQL	Result	Qual
10	Acetone	10	7	B	10	3	B			
10	Benzene	10	10	U	10	10	U			
10	Bromodichloromethane	10	10	U	10	10	U			
10	Bromoform	10	10	U	10	10	U			
10	Bromomethane	10	14	J	10	10	U			
10	2-Butanone	10	10	U	10	10	U			
10	Carbon Disulfide	10	10	U	10	10	U			
10	Carbon Tetrachloride	10	10	U	10	10	U			
10	Chlorobenzene	10	10	U	10	10	U			
10	Chloroethane	10	10	U	10	10	U			
10	Chloroform	10	10	U	10	10	U			
10	Chloromethane	10	20	J	10	10	U			
10	Dibromochloromethane	10	10	U	10	10	U			
10	1,1-Dichloroethane	10	10	U	10	10	U			
10	1,2-Dichloroethane	10	10	U	10	10	U			
10	1,1-Dichloroethene	10	10	U	10	10	U			
10	Total 1,2-Dichloroethene	10	10	U	10	10	U			
10	1,2-Dichloropropane	10	10	U	10	10	U			
10	cis-1,3-Dichloropropene	10	10	U	10	10	U			
10	trans-1,3-Dichloropropene	10	10	U	10	10	U			
10	Ethylbenzene	10	10	U	10	10	U			
10	2-Hexanone	10	10	U	10	10	U			
10	4-Methyl-2-Pentanone	10	10	U	10	10	U			
10	Methylene Chloride	10	22	B	10	2	B			
10	Styrene	10	10	U	10	10	U			
10	1,1,2,2-Tetrachloroethane	10	10	U	10	10	U			
10	Tetrachloroethene	10	10	U	10	10	U			
10	Toluene	10	10	U	10	10	U			
10	1,1,1-Trichloroethane	10	10	U	10	10	U			
10	1,1,2-Trichloroethane	10	10	U	10	10	U			
10	Trichloroethene	10	10	U	10	10	U			
10	Vinyl Chloride	10	10	U	10	10	U			
10	Total Xylenes	10	10	U	10	10	U			

Notes:
 All units are ug/l
 Sample Quantitation Limit (SQL) is CROL multiplied by the dilution factor and adjusted for percent moisture as appropriate.
 The Qual column indicates the qualifier applied to the result following data validation (see below).

Qualifiers:
 A - Acceptable (Quantitative) Data
 B - Present in associated blanks (Semi-quantitative Data)
 J - Estimated (Semi-quantitative) Data
 K - Biased High (Semi-quantitative) Data
 L - Biased Low (Semi-quantitative) Data
 M - Tentative Identification
 N - Tentative Identification
 R - Unusable Data
 U - Not Detected/Quantitative Data
 UL - Not Detected/Biased Low

AR302774

SUMMARY OF CLP ORGANICS ANALYSES - PHASE 1B
BERKS LANDFILL REMEDIAL INVESTIGATION
BERKS COUNTY, PENNSYLVANIA

Matrix: Groundwater Well

Volatle Organics

CROL	Parameter	Sample Point GUC01/1B			Sample Point GUC02/1B			Sample Point GUC03D/1B			Sample Point GUC03D/RE		
		SQL	Result	Qual	SQL	Result	Qual	SQL	Result	Qual	SQL	Result	Qual
10	Acetone	10	10	UJ	10	10	UJ	67	67	U	100	100	UJ
10	Benzene	10	10	U	10	10	U	67	67	U	100	100	U
10	Bromodichloromethane	10	10	U	10	10	U	67	67	U	100	100	U
10	Bromoform	10	10	U	10	10	U	67	67	U	100	100	U
10	Bromomethane	10	10	U	10	10	U	67	67	U	100	100	U
10	2-Butanone	10	10	U	10	10	U	67	67	J	100	100	U
10	Carbon Disulfide	10	10	U	10	10	U	67	67	U	100	100	U
10	Carbon Tetrachloride	10	10	U	10	10	U	67	67	U	100	100	U
10	Chlorobenzene	10	10	U	10	10	U	67	67	U	100	100	U
10	Chloroethane	10	10	U	10	10	U	67	67	U	100	100	U
10	Chloroform	10	10	U	10	10	U	67	67	U	100	100	U
10	Chloromethane	10	10	U	10	10	U	67	67	U	100	100	U
10	Dibromochloromethane	10	10	U	10	10	U	67	67	U	100	100	U
10	1,1-Dichloroethane	10	10	U	10	10	U	67	67	J	100	100	U
10	1,2-Dichloroethane	10	10	U	10	10	U	67	67	J	13	13	J
10	1,1-Dichloroethene	10	10	U	10	10	U	67	67	J	1600	1600	J
10	Total 1,2-Dichloroethene	10	10	U	10	10	U	67	67	J	100	100	J
10	1,2-Dichloropropane	10	10	U	10	10	U	67	67	U	100	100	U
10	cis-1,3-Dichloropropene	10	10	U	10	10	U	67	67	U	100	100	U
10	trans-1,3-Dichloropropene	10	10	U	10	10	U	67	67	U	100	100	U
10	Ethylbenzene	10	10	U	10	10	U	67	67	U	100	100	U
10	2-Hexanone	10	10	U	10	10	U	67	67	U	100	100	U
10	4-Methyl-2-Pentanone	10	10	UJ	10	10	UJ	67	67	UJ	100	100	UJ
10	Methylene Chloride	10	10	U	10	10	U	67	67	U	100	100	U
10	Styrene	10	17	B	10	3	B	67	37	B	200	200	B
10	1,1,2,2-Tetrachloroethane	10	10	U	10	10	U	67	67	U	100	100	U
10	Tetrachloroethene	10	10	U	10	10	U	67	67	U	100	100	U
10	Toluene	10	10	U	10	10	U	67	67	U	100	100	U
10	1,1,1-Trichloroethane	10	10	U	10	10	U	67	67	U	100	100	U
10	1,1,2-Trichloroethane	10	10	U	10	10	U	67	67	U	100	100	U
10	Trichloroethene	10	10	U	10	10	U	67	67	U	100	100	U
10	Vinyl Chloride	10	10	U	10	10	U	67	67	A	650	650	A
10	Total Xylenes	10	10	U	10	10	U	67	100	A	150	150	A
10		10	10	U	10	10	U	67	67	U	100	100	U

Notes:
All units are ug/l
Sample Quantitation Limit (SQL) is CROL multiplied by the dilution factor and adjusted for percent moisture as appropriate.
The Qual column indicates the qualifier applied to the result following data validation (see below).

Qualifiers:
A - Acceptable (Quantitative) Data
K - Biased High (Semi-quantitative) Data
UJ - Not Detected/Estimated Data
M - Tentative Identification
J - Estimated (Semi-quantitative) Data
L - Biased Low (Semi-quantitative) Data
U - Not Detected/Quantitative Data
R - Unusable Data
B - Present in associated blanks (Semi-quantitative Data)
UL - Not Detected/Biased Low
Q - Not Applicable

AR302775

SUMMARY OF CLP ORGANICS ANALYSES - PHASE 1B
BERKS LANDFILL REMEDIAL INVESTIGATION
BERKS COUNTY, PENNSYLVANIA

Matrix: Groundwater Well

Volatile Organics

CRQL	Parameter	Sample Point GVC03D/FD Date Sampled: 07/06/93 Dilution Factor: 1.0 Percent Moisture: NA %			Sample Point GVC03D/FR Date Sampled: 07/06/93 Dilution Factor: 10.0 Percent Moisture: NA %			Sample Point GVC03S/1B Date Sampled: 07/02/93 Dilution Factor: 1.0 Percent Moisture: NA %			Sample Point GVC04B/1B Date Sampled: 06/30/93 Dilution Factor: 1.0 Percent Moisture: NA %		
		SQL	Result	Qual	SQL	Result	Qual	SQL	Result	Qual	SQL	Result	Qual
10	Acetone	10	3	J	100	100	UJ	10	10	U	10	10	U
10	Benzene	10	10	U	100	100	U	10	2	J	10	10	U
10	Bromodichloromethane	10	10	U	100	100	U	10	10	U	10	10	U
10	Bromoform	10	10	U	100	100	U	10	10	U	10	10	U
10	Bromomethane	10	10	U	100	100	U	10	10	U	10	10	U
10	2-Butanone	10	10	U	100	100	U	10	10	U	10	10	U
10	Carbon Disulfide	10	10	U	100	100	U	10	10	U	10	10	U
10	Carbon Tetrachloride	10	10	U	100	100	U	10	10	U	10	10	U
10	Chlorobenzene	10	10	U	100	100	U	10	10	U	10	10	U
10	Chloroethane	10	6	J	100	100	U	10	5	J	10	10	U
10	Chloroform	10	10	U	100	100	U	10	4	J	10	10	U
10	Chloromethane	10	10	U	100	100	U	10	10	U	10	10	U
10	Dibromochloromethane	10	10	U	100	100	U	10	10	U	10	10	U
10	1,1-Dichloroethane	10	7	J	100	100	U	10	8	J	10	10	U
10	1,2-Dichloroethane	10	10	U	100	100	U	10	10	U	10	10	U
10	1,1-Dichloroethene	10	11	J	100	14	J	10	10	U	10	10	U
10	Total 1,2-Dichloroethene	10	1400	J	100	1800	A	10	53	A	10	10	U
10	1,2-Dichloropropane	10	10	U	100	100	U	10	10	U	10	10	U
10	cis-1,3-Dichloropropene	10	10	U	100	100	U	10	10	U	10	10	U
10	trans-1,3-Dichloropropene	10	10	U	100	100	U	10	10	U	10	10	U
10	Ethylbenzene	10	10	U	100	100	UJ	10	10	UJ	10	10	UJ
10	2-Hexanone	10	10	UJ	100	100	U	10	10	U	10	10	U
10	4-Methyl-2-Pentanone	10	10	U	100	100	U	10	10	U	10	10	U
10	Methylene Chloride	10	11	B	100	110	B	10	3	B	10	5	B
10	Styrene	10	10	U	100	100	U	10	10	U	10	10	U
10	1,1,2,2-Tetrachloroethane	10	1	J	100	100	U	10	10	U	10	10	U
10	Tetrachloroethene	10	10	U	100	100	U	10	10	U	10	10	U
10	Toluene	10	6	J	100	100	U	10	10	U	10	10	U
10	1,1,1-Trichloroethane	10	1	J	100	100	U	10	2	J	10	10	U
10	1,1,2-Trichloroethane	10	1	J	100	100	U	10	10	U	10	10	U
10	Trichloroethene	10	660	J	100	660	A	10	2	J	10	10	U
10	Vinyl Chloride	10	95	A	100	180	A	10	10	A	10	10	U
10	Total Xylenes	10	10	U	100	100	U	10	10	U	10	10	U

Notes:
 All units are ug/l
 Sample Quantitation Limit (SQL) is CRQL multiplied by the dilution factor and adjusted for percent moisture as appropriate.
 The Qual column indicates the qualifier applied to the result following data validation (see below).

Qualifiers:
 A - Acceptable (Quantitative) Data
 K - Biased High (Semi-quantitative) Data
 L - Biased Low (Semi-quantitative) Data
 J - Estimated (Semi-quantitative) Data
 B - Present in associated blanks (Semi-quantitative Data)
 UJ - Not Detected/Estimated Data
 U - Not Detected/Quantitative Data
 UL - Not Detected/Biased Low
 N - Tentative Identification
 R - Unusable Data
 Q - Not Applicable

AR302776

SUMMARY OF CLP ORGANICS ANALYSES - PHASE 1B
BERKS LANDFILL REMEDIAL INVESTIGATION
BERKS COUNTY, PENNSYLVANIA

Matrix: Groundwater Well

Volatile Organics

CRQL	Parameter	Sample Point GUC04S/1B			Sample Point GUC06b/1B			Sample Point GUC06S/1B			Sample Point GUC07b/1B		
		SQL	Result	Qual	SQL	Result	Qual	SQL	Result	Qual	SQL	Result	Qual
10	Acetone	10	10	U	10	10	U	10	10	U	10	10	U
10	Benzene	10	10	U	10	10	U	10	10	U	10	10	U
10	Bromodichloromethane	10	10	U	10	10	U	10	10	U	10	10	U
10	Bromoform	10	10	U	10	10	U	10	10	U	10	10	U
10	Bromomethane	10	10	U	10	10	U	10	10	U	10	10	U
10	2-Butanone	10	10	U	10	10	U	10	10	U	10	10	U
10	Carbon Disulfide	10	10	U	10	10	U	10	10	U	10	10	U
10	Carbon Tetrachloride	10	10	U	10	10	U	10	10	U	10	10	U
10	Chlorobenzene	10	10	U	10	10	U	10	10	U	10	10	U
10	Chloroethane	10	10	U	10	10	U	10	10	U	10	10	U
10	Chloroform	10	10	U	10	10	U	10	10	U	10	10	U
10	Chloromethane	10	10	U	10	10	U	10	10	U	10	10	U
10	Dibromochloromethane	10	10	U	10	10	U	10	10	U	10	10	U
10	1,1-Dichloroethane	10	1	J	10	10	U	10	10	U	10	10	U
10	1,2-Dichloroethane	10	10	U	10	10	U	10	10	U	10	10	U
10	1,1-Dichloroethene	10	8	J	10	10	U	10	10	U	10	10	U
10	Total 1,2-Dichloroethane	10	10	U	10	10	U	10	10	U	10	10	U
10	1,2-Dichloropropane	10	10	U	10	10	U	10	10	U	10	10	U
10	cis-1,3-Dichloropropene	10	10	U	10	10	U	10	10	U	10	10	U
10	trans-1,3-Dichloropropene	10	10	U	10	10	U	10	10	U	10	10	U
10	Ethylbenzene	10	10	U	10	10	U	10	10	U	10	10	U
10	2-Hexanone	10	10	U	10	10	U	10	10	U	10	10	U
10	4-Methyl-2-Pentanone	10	10	U	10	10	U	10	10	U	10	10	U
10	Methylene Chloride	10	4	U	10	10	U	10	10	U	10	10	U
10	Styrene	10	10	U	10	85	B	10	10	U	10	10	U
10	1,1,2,2-Tetrachloroethane	10	10	U	10	10	U	10	10	U	10	10	U
10	Tetrachloroethene	10	10	U	10	10	U	10	10	U	10	10	U
10	Toluene	10	10	U	10	10	U	10	10	U	10	10	U
10	1,1,1-Trichloroethane	10	10	U	10	10	U	10	10	U	10	10	U
10	1,1,2-Trichloroethane	10	10	U	10	10	U	10	10	U	10	10	U
10	Trichloroethene	10	2	J	10	10	U	10	10	U	10	10	U
10	Vinyl Chloride	10	2	J	10	10	U	10	10	U	10	10	U
10	Total Xylenes	10	10	U	10	10	U	10	10	U	10	10	U

Notes:

All units are ug/l
Sample Quantitation Limit (SQL) is CRQL multiplied by the dilution factor and adjusted for percent moisture as appropriate.
The Qual column indicates the qualifier applied to the result following data validation (see below).

Qualifiers:
A - Acceptable (Quantitative) Data K - Biased High (Semi-quantitative) Data UJ - Not Detected/Estimated Data M - Tentative Identification
J - Estimated (Semi-quantitative) Data L - Biased Low (Semi-quantitative) Data U - Not Detected/Quantitative Data R - Unusable Data
B - Present in associated blanks (Semi-quantitative Data) UL - Not Detected/Biased Low Q - Not Applicable

AR302777

SUMMARY OF CLP ORGANICS ANALYSES - PHASE 1B
BERKS LANDFILL REMEDIAL INVESTIGATION
BERKS COUNTY, PENNSYLVANIA

Matrix: Groundwater Well

Volatile Organics

CROL	Parameter	Sample Point G4C07S/1B			Sample Point G4C02/1B			Sample Point G4C01/1B			Sample Point G4C02/1B		
		SQL	Result	Qual	SQL	Result	Qual	SQL	Result	Qual	SQL	Result	Qual
10	Acetone	10	10	UJ	10	10	U	10	6	J	10	10	U
10	Benzene	10	10	U	10	10	J	10	10	U	10	10	U
10	Bromodichloromethane	10	10	U	10	10	U	10	10	U	10	10	U
10	Bromoform	10	10	U	10	10	U	10	10	U	10	10	U
10	Bromomethane	10	10	U	10	10	U	10	10	U	10	10	U
10	2-Butanone	10	10	U	10	10	U	10	10	U	10	10	U
10	Carbon Disulfide	10	10	U	10	10	U	10	10	U	10	10	U
10	Carbon Tetrachloride	10	10	U	10	10	U	10	10	U	10	10	U
10	Chlorobenzene	10	7	J	10	10	J	10	10	U	10	10	U
10	Chloroethane	10	10	U	10	10	U	10	10	U	10	10	U
10	Chloroform	10	10	U	10	10	U	10	10	U	10	10	U
10	Chloromethane	10	10	U	10	10	U	10	10	U	10	10	U
10	Dibromochloromethane	10	10	U	10	10	U	10	10	U	10	10	U
10	1,1-Dichloroethane	10	10	U	10	10	U	10	10	U	10	10	U
10	1,2-Dichloroethane	10	10	U	10	10	U	10	10	U	10	10	U
10	1,1-Dichloroethene	10	10	U	10	10	U	10	10	U	10	10	U
10	Total 1,2-Dichloroethene	10	10	U	10	3	J	10	10	U	10	10	U
10	1,2-Dichloropropane	10	1	J	10	10	U	10	10	U	10	10	U
10	cis-1,3-Dichloropropene	10	10	U	10	10	U	10	10	U	10	10	U
10	trans-1,3-Dichloropropene	10	10	U	10	10	U	10	10	U	10	10	U
10	Ethylbenzene	10	10	U	10	10	U	10	10	U	10	10	U
10	2-Hexanone	10	10	UJ	10	10	UJ	10	10	UJ	10	10	UJ
10	4-Methyl-2-Pentanone	10	10	U	10	10	U	10	10	U	10	10	U
10	Methylene Chloride	10	4	B	10	10	B	10	7	B	10	10	B
10	Styrene	10	10	U	10	10	U	10	10	U	10	10	U
10	1,1,2,2-Tetrachloroethane	10	10	U	10	10	U	10	10	U	10	10	U
10	Tetrachloroethene	10	10	U	10	10	U	10	10	U	10	10	U
10	Toluene	10	10	U	10	10	U	10	10	U	10	10	U
10	1,1,1-Trichloroethane	10	10	U	10	10	U	10	10	U	10	10	U
10	1,1,2-Trichloroethane	10	10	U	10	10	U	10	10	U	10	10	U
10	Trichloroethene	10	10	U	10	10	U	10	10	U	10	10	U
10	Vinyl Chloride	10	10	U	10	10	U	10	10	U	10	10	U
10	Total Xylenes	10	10	U	10	10	U	10	10	U	10	10	U

Notes:

All units are ug/l
Sample Quantitation Limit (SQL) is CROL multiplied by the dilution factor and adjusted for percent moisture as appropriate.
The Qual column indicates the qualifier applied to the result following data validation (see below).

Qualifiers:
A - Accepted (Quantitative) Data
K - Biased High (Semi-quantitative) Data
L - Biased Low (Semi-quantitative) Data
B - Present in associated blanks (Semi-quantitative Data)

UJ - Not Detected/Estimated Data
U - Not Detected/quantitative Data
UL - Not Detected/Biased Low
M - Tentative Identification
N - Unusable Data
R - Not Applicable
Q - Not Applicable

AR302778

Matrix: Groundwater Well

Volatile Organics

CRQL	Parameter	Sample Point GUG03/18			Sample Point GUG04/18			Sample Point GUG05/18			Sample Point GUG06/18		
		SQL	Result	Qual	SQL	Result	Qual	SQL	Result	Qual	SQL	Result	Qual
10	Acetone	10	10	U	10	10	U	10	10	U	10	10	U
10	Benzene	10	10	U	10	10	U	10	10	U	10	10	U
10	Bromodichloromethane	10	10	U	10	10	U	10	10	U	10	10	U
10	Bromoform	10	10	U	10	10	U	10	10	U	10	10	U
10	Bromoethane	10	10	U	10	10	U	10	10	U	10	10	U
10	2-Butenone	10	10	U	10	10	U	10	10	U	10	10	U
10	Carbon Disulfide	10	10	U	10	10	U	10	10	U	10	10	U
10	Carbon Tetrachloride	10	10	U	10	10	U	10	10	U	10	10	U
10	Chlorobenzene	10	10	U	10	10	U	10	10	U	10	10	U
10	Chloroethane	10	10	U	10	10	U	10	10	U	10	10	U
10	Chloroform	10	10	U	10	10	U	10	10	U	10	10	U
10	Chloroethane	10	10	U	10	10	U	10	10	U	10	10	U
10	Dibromochloromethane	10	10	U	10	10	U	10	10	U	10	10	U
10	1,1-Dichloroethane	10	10	U	10	10	U	10	10	U	10	10	U
10	1,2-Dichloroethane	10	10	U	10	10	U	10	10	U	10	10	U
10	1,1-Dichloroethene	10	10	U	10	10	U	10	10	U	10	10	U
10	Total 1,2-Dichloroethene	10	10	U	10	10	U	10	10	U	10	10	U
10	1,2-Dichloropropane	10	10	U	10	10	U	10	10	U	10	10	U
10	cis-1,3-Dichloropropene	10	10	U	10	10	U	10	10	U	10	10	U
10	trans-1,3-Dichloropropene	10	10	U	10	10	U	10	10	U	10	10	U
10	Ethylbenzene	10	10	U	10	10	U	10	10	U	10	10	U
10	2-Hexanone	10	10	U	10	10	U	10	10	U	10	10	U
10	4-Methyl-2-Pentanone	10	10	U	10	10	U	10	10	U	10	10	U
10	Methylene Chloride	10	10	U	10	10	U	10	10	U	10	10	U
10	Styrene	10	10	U	10	10	U	10	10	U	10	10	U
10	1,1,2-Tetrachloroethane	10	10	U	10	10	U	10	10	U	10	10	U
10	Tetrachloroethene	10	10	U	10	10	U	10	10	U	10	10	U
10	Toluene	10	10	U	10	10	U	10	10	U	10	10	U
10	1,1,1-Trichloroethane	10	10	U	10	10	U	10	10	U	10	10	U
10	1,1,2-Trichloroethane	10	10	U	10	10	U	10	10	U	10	10	U
10	Trichloroethene	10	10	U	10	10	U	10	10	U	10	10	U
10	Vinyl Chloride	10	10	U	10	10	U	10	10	U	10	10	U
10	Total Xylenes	10	10	U	10	10	U	10	10	U	10	10	U

Notes:
 All units are ug/l
 Sample Quantitation Limit (SQL) is CRQL multiplied by the dilution factor and adjusted for percent moisture as appropriate.
 The Qual column indicates the qualifier applied to the result following data validation (see below).

Qualifiers:
 A - Acceptable (Quantitative) Data
 J - Estimated (Semi-quantitative) Data
 B - Present in associated blanks (Semi-quantitative Data)
 K - Biased High (Semi-quantitative) Data
 L - Biased Low (Semi-quantitative) Data
 UJ - Not Detected/Estimated Data
 U - Not Detected/Quantitative Data
 UL - Not Detected/Biased Low
 N - Tentative Identification
 R - Unusable Data
 Q - Not Applicable

SUMMARY OF CLP OR CAS ANALYSES - PHASE 1B
BERKS LANDFILL REMEDIAL INVESTIGATION
BERKS COUNTY, PENNSYLVANIA

Matrix: Groundwater Well

Volatile Organics

CROL	Parameter	Sample Point GW07/1B			Sample Point GW08/1B			Sample Point GW10/1B			Sample Point GW11/1B		
		SQL	Result	Qual	SQL	Result	Qual	SQL	Result	Qual	SQL	Result	Qual
10	Acetone	10	10	U	10	10	U	10	10	U	10	10	U
10	Benzene	10	10	U	10	10	U	10	10	U	10	10	U
10	Bromodichloromethane	10	10	U	10	10	U	10	10	U	10	10	U
10	Bromoform	10	10	U	10	10	U	10	10	U	10	10	U
10	Bromomethane	10	10	U	10	10	U	10	10	U	10	10	U
10	2-Butanone	10	10	U	10	10	U	10	10	U	10	10	U
10	Carbon Disulfide	10	10	U	10	10	U	10	10	U	10	10	U
10	Carbon Tetrachloride	10	10	U	10	10	U	10	10	U	10	10	U
10	Chlorobenzene	10	10	U	10	10	U	10	10	U	10	10	U
10	Chloroethane	10	10	U	10	10	U	10	10	U	10	10	U
10	Chloroform	10	10	U	10	10	U	10	10	U	10	10	U
10	Chloromethane	10	10	U	10	10	U	10	10	U	10	10	U
10	Dibromochloromethane	10	10	U	10	10	U	10	10	U	10	10	U
10	1,1-Dichloroethane	10	10	U	10	10	U	10	10	U	10	10	U
10	1,2-Dichloroethane	10	10	U	10	10	U	10	10	U	10	10	U
10	Total 1,2-Dichloroethane	10	10	U	10	10	U	10	10	U	10	10	U
10	1,2-Dichloropropane	10	10	U	10	10	U	10	10	U	10	10	U
10	cis-1,3-Dichloropropene	10	10	U	10	10	U	10	10	U	10	10	U
10	trans-1,3-Dichloropropene	10	10	U	10	10	U	10	10	U	10	10	U
10	Ethylbenzene	10	10	U	10	10	U	10	10	U	10	10	U
10	2-Hexanone	10	10	U	10	10	U	10	10	U	10	10	U
10	4-Methyl-2-Pentanone	10	10	U	10	10	U	10	10	U	10	10	U
10	Methylene Chloride	10	10	U	10	10	U	10	10	U	10	10	U
10	Styrene	10	2	B	10	1	B	10	7	B	10	21	B
10	1,1,2,2-Tetrachloroethane	10	10	U	10	10	U	10	10	U	10	10	U
10	Tetrachloroethane	10	10	U	10	10	U	10	10	U	10	10	U
10	Toluene	10	10	U	10	10	U	10	10	U	10	10	U
10	1,1,1-Trichloroethane	10	10	U	10	10	U	10	10	U	10	10	U
10	1,1,2-Trichloroethane	10	10	U	10	10	U	10	10	U	10	10	U
10	Trichloroethene	10	10	U	10	10	U	10	10	U	10	10	U
10	Vinyl Chloride	10	10	U	10	10	U	10	10	U	10	10	U
10	Total Xylenes	10	10	U	10	10	U	10	10	U	10	10	U

Notes:

All units are ug/l
Sample Quantitation Limit (SQL) is CROL multiplied by the dilution factor and adjusted for percent moisture as appropriate.
The qual column indicates the qualifier applied to the result following data validation (see below).

Qualifiers:.....
A - Acceptable (Quantitative) Data K - Biased High (Semi-quantitative) Data UJ - Not Detected/Estimated Data M - Tentative Identification
J - Estimated (Semi-quantitative) Data L - Biased Low (Semi-quantitative) Data U - Not Detected/Quantitative Data R - Unusable Data
B - Present in associated blanks (Semi-quantitative Data) UL - Not Detected/Biased Low Q - Not Applicable

AR302780

SUMMARY OF CLP ORGANICS ANALYSES - PHASE 1B
BERKS LANDFILL REMEDIAL INVESTIGATION
BERKS COUNTY, PENNSYLVANIA

Matrix: Groundwater Well

Volatile Organics

CROL	Parameter	Sample Point GUG11/FD			Sample Point GUG12/FB			Sample Point GUG12/FB			Sample Point GUG13/FB		
		SQL	Result	Qual	SQL	Result	Qual	SQL	Result	Qual	SQL	Result	Qual
10	Acetone	10	10	U	10	10	U	10	10	U	10	10	U
10	Benzene	10	2	J	10	1	J	10	1	J	10	10	U
10	Bromodichloromethane	10	10	U	10	10	U	10	10	U	10	10	U
10	Bromoform	10	10	U	10	10	U	10	10	U	10	10	U
10	Bromomethane	10	10	U	10	10	U	10	10	U	10	10	U
10	2-Butanone	10	10	U	10	10	U	10	10	U	10	10	U
10	Carbon Disulfide	10	10	U	10	3	J	10	10	U	10	10	U
10	Carbon Tetrachloride	10	10	U	10	10	U	10	10	U	10	10	U
10	Chlorobenzene	10	22	A	10	3	J	10	3	J	10	10	U
10	Chloroethane	10	10	UJ	10	6	J	10	6	J	10	10	U
10	Chloroform	10	10	UJ	10	10	U	10	10	U	10	10	U
10	Chloromethane	10	10	U	10	10	U	10	10	U	10	10	U
10	Dibromochloromethane	10	10	U	10	10	U	10	10	U	10	10	U
10	1,1-Dichloroethane	10	10	U	10	10	U	10	10	U	10	10	U
10	1,2-Dichloroethane	10	10	U	10	10	U	10	10	U	10	10	U
10	Total 1,2-Dichloroethane	10	10	U	10	1	J	10	1	J	10	3	J
10	1,2-Dichloropropane	10	10	U	10	10	U	10	10	U	10	10	U
10	cis-1,3-Dichloropropene	10	10	U	10	10	U	10	10	U	10	10	U
10	trans-1,3-Dichloropropene	10	10	U	10	10	U	10	10	U	10	10	U
10	Ethylbenzene	10	10	U	10	10	U	10	10	U	10	10	U
10	2-Hexanone	10	10	U	10	10	U	10	10	U	10	10	U
10	4-Methyl-2-Pentanone	10	10	UJ	10	10	UJ	10	10	UJ	10	10	UJ
10	Methylene Chloride	10	12	B	10	3	B	10	3	B	10	14	B
10	Styrene	10	10	U	10	10	U	10	10	U	10	10	U
10	1,1,2,2-Tetrachloroethane	10	10	U	10	10	U	10	10	U	10	10	U
10	Tetrachloroethane	10	10	U	10	10	U	10	10	U	10	10	U
10	Toluene	10	10	U	10	10	U	10	10	U	10	10	U
10	1,1,1-Trichloroethane	10	10	U	10	10	U	10	10	U	10	10	U
10	1,1,2-Trichloroethane	10	10	U	10	10	U	10	10	U	10	10	U
10	Trichloroethene	10	10	U	10	10	U	10	10	U	10	10	U
10	Vinyl Chloride	10	10	U	10	10	U	10	10	U	10	15	U
10	Total Xylenes	10	10	U	10	10	U	10	10	U	10	10	U

Notes:
All units are ug/l
Sample Quantitation Limit (SQL) is CROL multiplied by the dilution factor and adjusted for percent moisture as appropriate.
The dual column indicates the qualifier applied to the result following data validation (see below).

Qualifiers:
A - Acceptable (Quantitative) Data
K - Biased High (Semi-quantitative) Data
UJ - Not Detected/Estimated Data
N - Tentative Identification
J - Estimated (Semi-quantitative) Data
L - Biased Low (Semi-quantitative) Data
U - Not Detected/quantitative Data
R - Unusable Data
B - Present in associated blanks (Semi-quantitative Data)
UL - Not Detected/Biased Low
Q - Not Applicable

SUMMARY OF CLP ORGANICS ANALYSES - PHASE 1B
BERKS LANDFILL REMEDIAL INVESTIGATION
BERKS COUNTY, PENNSYLVANIA

Matrix: Groundwater Well

Volatile Organics

CRQL	Parameter	Sample Point GM03/18			Sample Point GM06/18			Sample Point GM11/18			Sample Point GM140/18		
		SOL	Result	Qual	SOL	Result	Qual	SOL	Result	Qual	SOL	Result	Qual
10	Acetone	10	19	J	10	10	UJ	10	10	U	10	10	UJ
10	Benzene	10	4	J	10	10	U	10	10	U	10	10	U
10	Bromodichloromethane	10	10	U	10	10	U	10	10	U	10	10	U
10	Bromoform	10	10	U	10	10	U	10	10	U	10	10	U
10	Bromomethane	10	10	U	10	10	U	10	10	U	10	10	U
10	2-Butanone	10	10	U	10	10	U	10	10	U	10	10	U
10	Carbon Disulfide	10	10	U	10	10	U	10	10	U	10	10	U
10	Carbon Tetrachloride	10	10	U	10	10	U	10	10	U	10	10	U
10	Chlorobenzene	10	4	U	10	10	U	10	10	U	10	10	U
10	Chloroethane	10	27	J	10	10	U	10	10	U	10	10	J
10	Chloroform	10	10	A	10	10	U	10	10	U	10	10	U
10	Chloromethane	10	10	U	10	10	U	10	10	U	10	10	U
10	Dibromochloromethane	10	10	U	10	10	U	10	10	U	10	10	U
10	1,1-Dichloroethane	10	9	J	10	10	U	10	10	J	10	10	U
10	1,2-Dichloroethane	10	10	U	10	10	U	10	10	U	10	10	U
10	1,1,1-Trichloroethane	10	53	A	10	10	U	10	10	A	10	10	U
10	Total 1,2-Dichloroethane	10	10	U	10	10	U	10	10	U	10	10	U
10	1,2-Dichloropropene	10	10	U	10	10	U	10	10	U	10	10	U
10	cis-1,3-Dichloropropene	10	10	U	10	10	U	10	10	U	10	10	U
10	trans-1,3-Dichloropropene	10	10	U	10	10	U	10	10	U	10	10	U
10	Ethylbenzene	10	10	U	10	10	U	10	10	U	10	10	U
10	2-Hexanone	10	10	UJ	10	10	UJ	10	10	UJ	10	10	UJ
10	4-Methyl-2-Pentanone	10	10	U	10	10	U	10	10	U	10	10	U
10	Methylene Chloride	10	4	B	10	7	B	10	3	B	10	5	B
10	Styrene	10	10	U	10	10	U	10	10	U	10	10	U
10	1,1,2-Tetrachloroethane	10	10	U	10	10	U	10	10	U	10	10	U
10	Tetrachloroethane	10	10	U	10	10	U	10	10	U	10	10	U
10	Toluene	10	10	U	10	10	U	10	10	U	10	10	U
10	1,1,1-Trichloroethane	10	10	U	10	10	U	10	10	U	10	10	U
10	1,1,2-Trichloroethane	10	10	U	10	10	U	10	10	U	10	10	U
10	Trichloroethene	10	61	U	10	10	U	10	10	U	10	10	U
10	Vinyl Chloride	10	9	J	10	10	U	10	10	A	10	10	U
10	Total Xylenes	10	10	U	10	10	U	10	10	J	10	10	U

Notes:
All units are ug/l
Sample Quantitation Limit (SOL) is CRQL multiplied by the dilution factor and adjusted for percent moisture as appropriate.
The Qual column indicates the qualifier applied to the result following data validation (see below).

Qualifiers:
A - Acceptable (Quantitative) Data
K - Biased High (Semi-quantitative) Data
UJ - Not Detected/Estimated Data
N - Tentative Identification
J - Estimated (Semi-quantitative) Data
L - Biased Low (Semi-quantitative) Data
U - Not Detected/Quantitative Data
R - Unusable Data
B - Present in associated blanks (Semi-quantitative Data)
UL - Not Detected/Biased Low
Q - Not Applicable

Matrix: Groundwater Well

Volatile Organics

CRQL	Parameter	Sample Point GM14S/1B Date Sampled: 07/09/93 Dilution Factor: 1.0 Percent Moisture: NA %			Sample Point GM150/1B Date Sampled: 07/08/93 Dilution Factor: 1.0 Percent Moisture: NA %			Sample Point GM15S/1B Date Sampled: 07/02/93 Dilution Factor: 1.0 Percent Moisture: NA %			Sample Point GM16/1B Date Sampled: 07/08/93 Dilution Factor: 1.0 Percent Moisture: NA %		
		SQL	Result	Qual	SQL	Result	Qual	SQL	Result	Qual	SQL	Result	Qual
10	Acetone	10	10	UJ	10	10	UJ	10	10	UJ	10	10	UJ
10	Benzene	10	5	J	10	10	U	10	10	U	10	11	A
10	Bromochloromethane	10	10	U	10	10	U	10	10	U	10	10	U
10	Bromoform	10	10	U	10	10	U	10	10	U	10	10	U
10	Bromomethane	10	10	U	10	10	U	10	10	U	10	10	U
10	2-Butanone	10	10	U	10	10	U	10	10	U	10	10	U
10	Carbon Disulfide	10	10	U	10	10	U	10	10	U	10	10	U
10	Carbon Tetrachloride	10	22	A	10	10	U	10	10	U	10	3	J
10	Chlorobenzene	10	10	U	10	10	U	10	10	U	10	10	U
10	Chloroethane	10	10	U	10	10	U	10	10	U	10	10	U
10	Chloroform	10	10	U	10	10	U	10	10	U	10	10	U
10	Chloromethane	10	10	U	10	10	U	10	10	U	10	10	U
10	Dibromochloromethane	10	10	U	10	10	U	10	10	U	10	10	U
10	1,1-Dichloroethane	10	10	U	10	10	U	10	10	U	10	10	U
10	1,2-Dichloroethane	10	10	U	10	10	U	10	10	U	10	10	U
10	1,1-Dichloroethene	10	10	U	10	10	U	10	10	U	10	10	U
10	Total 1,2-Dichloroethene	10	10	U	10	10	U	10	10	U	10	10	U
10	1,2-Dichloropropane	10	10	U	10	10	U	10	10	U	10	10	U
10	cis-1,3-Dichloropropene	10	10	U	10	10	U	10	10	U	10	10	U
10	trans-1,3-Dichloropropene	10	10	U	10	10	U	10	10	U	10	10	U
10	Ethylbenzene	10	10	UJ	10	10	UJ	10	10	UJ	10	5	J
10	2-Hexanone	10	10	U	10	10	U	10	10	U	10	10	UJ
10	4-Methyl-2-Pentanone	10	6	U	10	10	U	10	10	U	10	10	U
10	Methylene Chloride	10	10	B	10	10	B	10	10	B	10	4	B
10	Styrene	10	10	U	10	10	U	10	10	U	10	10	U
10	1,1,2,2-Tetrachloroethane	10	10	U	10	10	U	10	10	U	10	10	U
10	1,1,2-Trichloroethane	10	10	U	10	10	U	10	10	U	10	10	U
10	Toluene	10	10	U	10	10	U	10	10	U	10	10	U
10	1,1,1-Trichloroethane	10	10	U	10	10	U	10	10	U	10	10	U
10	1,1,2-Trichloroethane	10	10	U	10	10	U	10	10	U	10	10	U
10	Trichloroethane	10	10	U	10	10	U	10	10	U	10	10	U
10	Vinyl Chloride	10	10	U	10	10	U	10	10	U	10	10	U
10	Total Xylenes	10	10	U	10	10	U	10	10	U	10	9%	A

Notes:
All units are ug/l
Sample Quantitation Limit (SQL) is CRQL multiplied by the dilution factor and adjusted for percent moisture as appropriate.
The Qual column indicates the qualifier applied to the result following data validation (see below).

Qualifiers:
A - Acceptable (Quantitative) Data
K - Biased High (Semi-quantitative) Data
J - Estimated (Semi-quantitative) Data
L - Biased Low (Semi-quantitative) Data
B - Present in associated blanks (Semi-quantitative Data)
UJ - Not Detected/Estimated Data
U - Not Detected/Quantitative Data
UL - Not Detected/Biased Low
N - Tentative Identification
R - Unusable Data
Q - Not Applicable

AR302783

SUMMARY OF CLP ORGANICS ANALYSES - PHASE 1B
BERKS LANDFILL REMEDIAL INVESTIGATION
BERKS COUNTY, PENNSYLVANIA

Matrix: Groundwater Well

Volatile Organics

CRQL	Parameter	Sample Point GM177/1B Date Sampled: 07/08/93 Dilution Factor: 1.0 Percent Moisture: NA %		Sample Point GM195/1B Date Sampled: 06/30/93 Dilution Factor: 1.0 Percent Moisture: NA %		Sample Point GM190/1B Date Sampled: 07/06/93 Dilution Factor: 1.0 Percent Moisture: NA %		Sample Point Date Sampled: Dilution Factor: Percent Moisture: %			
		SQL	Result	SQL	Result	SQL	Result	SQL	Result	Qual	
10	Acetone	10	10	10	10	10	10	10	10	U	Qual
10	Benzene	10	10	10	10	10	10	10	10	U	Qual
10	Bromodichloromethane	10	10	10	10	10	10	10	10	U	Qual
10	Bromoform	10	10	10	10	10	10	10	10	U	Qual
10	Bromomethane	10	10	10	10	10	10	10	10	U	Qual
10	2-Butenone	10	10	10	10	10	10	10	10	U	Qual
10	Carbon Disulfide	10	10	10	10	10	10	10	10	U	Qual
10	Carbon Tetrachloride	10	10	10	10	10	10	10	10	U	Qual
10	Chlorobenzene	10	10	10	10	10	10	10	10	U	Qual
10	Chloroethane	10	10	10	10	10	10	10	10	U	Qual
10	Chloroform	10	10	10	10	10	10	10	10	U	Qual
10	Chloromethane	10	10	10	10	10	10	10	10	U	Qual
10	Dibromochloromethane	10	10	10	10	10	10	10	10	U	Qual
10	1,1-Dichloroethane	10	10	10	10	10	10	10	10	U	Qual
10	1,2-Dichloroethane	10	10	10	10	10	10	10	10	U	Qual
10	1,1-Dichloroethene	10	10	10	10	10	10	10	10	U	Qual
10	Total 1,2-Dichloroethene	10	10	10	10	10	10	10	10	U	Qual
10	1,2-Dichloropropane	10	10	10	10	10	10	10	10	U	Qual
10	cis-1,3-Dichloropropene	10	10	10	10	10	10	10	10	U	Qual
10	trans-1,3-Dichloropropene	10	10	10	10	10	10	10	10	U	Qual
10	Ethylbenzene	10	10	10	10	10	10	10	10	U	Qual
10	2-Hexenone	10	10	10	10	10	10	10	10	U	Qual
10	4-Methyl-2-Pentanone	10	10	10	10	10	10	10	10	U	Qual
10	Methylene Chloride	10	5	10	2	10	3	10	10	U	Qual
10	Styrene	10	10	10	10	10	10	10	10	U	Qual
10	1,1,2,2-Tetrachloroethane	10	10	10	10	10	10	10	10	U	Qual
10	Tetrachloroethene	10	10	10	10	10	10	10	10	U	Qual
10	Toluene	10	10	10	10	10	10	10	10	U	Qual
10	1,1,1-Trichloroethane	10	10	10	10	10	10	10	10	U	Qual
10	1,1,2-Trichloroethane	10	10	10	10	10	10	10	10	U	Qual
10	Trichloroethene	10	10	10	10	10	10	10	10	U	Qual
10	Vinyl Chloride	10	10	10	10	10	10	10	10	U	Qual
10	Total Xylenes	10	10	10	10	10	10	10	10	U	Qual

Notes:

All units are ug/l
Sample Quantitation Limit (SQL) is CRQL multiplied by the dilution factor and adjusted for percent moisture as appropriate.
The Qual column indicates the qualifier applied to the result following data validation (see below).

Qualifiers:
A - Acceptable (Quantitative) Data
J - Estimated (Semi-quantitative) Data
B - Present in associated blanks (Semi-quantitative Data)
K - Biased High (Semi-quantitative) Data
L - Biased Low (Semi-quantitative) Data
UJ - Not Detected/Estimated Data
U - Not Detected/Quantitative Data
UL - Not Detected/Biased Low
N - Tentative Identification
R - Unusable Data
Q - Not Applicable

AR302784

SEMI-VOLATILE ORGANIC COMPOUNDS

Phase 1A
Phase 1B

AR302785

SUMMARY OF CLP ORGANIC ANALYSES - PHASE 1A
BERKS LANDFILL REMEDIAL INVESTIGATION
BERKS COUNTY, PENNSYLVANIA

Matrix: Groundwater Well

Semivolatile Organics

CRQL	Parameter	Sample Point GR180/1A Date Sampled: 11/13/92 Dilution Factor: Percent Moisture: NA %			Sample Point GM-16/1A Date Sampled: 09/04/92 Dilution Factor: Percent Moisture: NA %			Sample Point GM-16/1A Date Sampled: 09/09/92 Dilution Factor: Percent Moisture: NA %			Sample Point GM-C5/1A Date Sampled: 09/03/92 Dilution Factor: Percent Moisture: NA %		
		SQL	Result	Qual	SQL	Result	Qual	SQL	Result	Qual	SQL	Result	Qual
10	Acenaphthene	10	10	U	10	10	U	10	10	U	10	10	U
10	Acenaphthylene	10	10	U	10	10	U	10	10	U	10	10	U
10	Anthracene	10	10	U	10	10	U	10	10	U	10	10	U
10	Benzo(a)Anthracene	10	10	U	10	10	U	10	10	U	10	10	U
10	4-Bromophenyl-phthalate	10	10	U	10	10	U	10	10	U	10	10	U
10	Butylbenzylphthalate	10	10	U	10	10	U	10	10	U	10	10	U
10	Carbazole	10	10	U	10	10	U	10	10	U	10	10	U
10	4-Chloro-3-Methylphenol	10	10	U	10	10	U	10	10	U	10	10	U
10	4-Chloroaniline	10	10	U	10	10	U	10	10	U	10	10	U
10	bis(2-Chloroethoxy)Methane	10	10	U	10	10	U	10	10	U	10	10	U
10	bis(2-Chloroethyl)ether	10	10	U	10	10	U	10	10	U	10	10	U
10	2-Chloronaphthalene	10	10	U	10	10	U	10	10	U	10	10	U
10	2-Chlorophenol	10	10	U	10	10	U	10	10	U	10	10	U
10	4-Chlorophenyl-phenylether	10	10	U	10	10	U	10	10	U	10	10	U
10	Chrysene	10	10	U	10	10	U	10	10	U	10	10	U
10	Dibenz(a,h)Anthracene	10	10	U	10	10	U	10	10	U	10	10	U
10	Dibenzofuran	10	10	U	10	10	U	10	10	U	10	10	U
10	1,2-Dichlorobenzene	10	10	U	10	10	U	10	10	U	10	10	U
10	1,3-Dichlorobenzene	10	10	U	10	10	U	10	10	U	10	10	U
10	1,4-Dichlorobenzene	10	10	U	10	10	U	10	10	U	10	10	U
10	3,3'-Dichlorobenzidine	10	10	U	10	10	U	10	10	U	10	10	U
10	2,4-Dichlorophenol	10	10	U	10	10	U	10	10	U	10	10	U
10	Diethylphthalate	10	10	U	10	10	U	10	10	U	10	10	U
10	2,4-Dimethylphenol	10	10	U	10	10	U	10	10	U	10	10	U
10	Dimethyl Phthalate	10	10	U	10	10	U	10	10	U	10	10	U
10	Di-n-butylphthalate	10	10	U	10	10	U	10	10	U	10	10	U
25	4,6-Dinitro-2-methylphenol	25	25	U	25	25	U	25	25	U	25	25	U
25	2,4-Dinitrophenol	25	25	U	25	25	U	25	25	U	25	25	U
10	2,4-Dinitrotoluene	10	10	U	10	10	U	10	10	U	10	10	U
10	2,6-Dinitrotoluene	10	10	U	10	10	U	10	10	U	10	10	U
10	Di-n-Octyl Phthalate	10	10	U	10	10	U	10	10	U	10	10	U
10	bis(2-Ethylhexyl)Phthalate	10	10	U	10	10	U	10	10	U	10	10	U
10	Fluoranthene	10	10	U	10	10	U	10	10	U	10	10	U
10	Benzo(b)Fluoranthene	10	10	U	10	10	U	10	10	U	10	10	U
10	Benzo(k)Fluoranthene	10	10	U	10	10	U	10	10	U	10	10	U
10	Fluorene	10	10	U	10	10	U	10	10	U	10	10	U
10	Hexachlorobenzene	10	10	U	10	10	U	10	10	U	10	10	U
10	Hexachlorobutadiene	10	10	U	10	10	U	10	10	U	10	10	U
10	Hexachlorocyclopentadiene	10	10	U	10	10	U	10	10	U	10	10	U
10	Hexachloroethane	10	10	U	10	10	U	10	10	U	10	10	U

Notes:

All units are ug/l
Sample Quantitation Limit (SQL) is CRQL multiplied by the dilution factor and adjusted for percent moisture as appropriate.
The Qual column indicates the qualifier following data validation (see below).

Qualifiers:.....
A - Acceptable (Quantitative) Data
J - Estimated (Semi-quantitative) Data
B - Present in associated blanks (Semi-quantitative Data)
K - Biased High (Semi-quantitative) Data
L - Biased Low (Semi-quantitative) Data
UL - Not Detected/Biased Low
UJ - Not Detected/Estimated Data
U - Not Detected/Quantitative Data
UL - Not Detected/Biased Low
M - Tentative Identification
N - Unusable Data
Q - Not Applicable

AR302786

Matrix: Groundwater Well

Semi-volatile Organics

CROL	Parameter	Sample Point GR18D/1A			Sample Point GU-16/1A			Sample Point GU-16/1A			Sample Point GU-C5/1A		
		Date Sampled: 11/13/92	Dilution Factor:	Percent Moisture: MA %	Date Sampled: 09/04/92	Dilution Factor:	Percent Moisture: MA %	Date Sampled: 09/09/92	Dilution Factor:	Percent Moisture: MA %	Date Sampled: 09/03/92	Dilution Factor:	Percent Moisture: MA %
		SQL	Result	Qual	SQL	Result	Qual	SQL	Result	Qual	SQL	Result	Qual
10	Indeno(1,2,3-cd)Pyrene	10	10	U	10	10	U	10	10	U	10	10	U
10	Isophorone	10	10	U	10	2	J	10	10	J	10	10	U
10	2-Methylnaphthalene	10	10	U	10	10	U	10	10	U	10	10	U
10	2-Methylphenol	10	10	U	10	10	U	10	10	U	10	10	U
10	4-Methylphenol	10	10	U	10	10	U	10	10	U	10	10	U
10	Naphthalene	10	10	U	10	2	J	10	10	U	10	10	U
25	2-Nitroaniline	25	25	U	25	25	U	25	25	U	25	25	U
25	3-Nitroaniline	25	25	U	25	25	U	25	25	U	25	25	U
25	4-Nitroaniline	25	25	U	25	25	U	25	25	U	25	25	U
10	Nitrobenzene	10	10	U	10	10	U	10	10	U	10	10	U
10	2-Nitrophenol	10	10	U	10	10	U	10	10	U	10	10	U
25	4-Nitrophenol	25	25	U	25	25	U	25	25	U	25	25	U
10	N-Nitroso-di-n-propylamine	10	10	U	10	10	U	10	10	U	10	10	U
10	N-Nitrosodiphenylamine	10	10	U	10	10	U	10	10	U	10	10	U
10	2,2'-Oxybis(1-Chloropropane)	10	10	U	10	10	U	10	10	U	10	10	U
10	Pentachlorophenol	25	25	U	25	25	U	25	25	U	25	25	U
10	Benzot(g,h,i)perylene	10	10	U	10	10	U	10	10	U	10	10	U
10	Phenanthrene	10	10	U	10	10	U	10	10	U	10	10	U
10	Phenol	10	10	U	10	10	U	10	10	U	10	10	U
10	Pyrene	10	10	U	10	10	U	10	10	U	10	10	U
10	Benzot(a)Pyrene	10	10	U	10	10	U	10	10	U	10	10	U
10	1,2,4-Trichlorobenzene	10	10	U	10	10	U	10	10	U	10	10	U
25	2,4,5-Trichlorophenol	25	25	U	25	25	U	25	25	U	25	25	U
10	2,4,6-Trichlorophenol	10	10	U	10	10	U	10	10	U	10	10	U

Notes:

All units are ug/l
Sample Quantitation Limit (SQL) is CROL multiplied by the dilution factor and adjusted for percent moisture as appropriate.
The Qual column indicates the qualifier applied to the result following data validation (see below).

Qualifiers:.....
A - Acceptable (Quantitative) Data K - Biased High (Semi-quantitative) Data UJ - Not Detected/Estimated Data M - Tentative Identification
J - Estimated (Semi-quantitative) Data L - Biased Low (Semi-quantitative) Data U - Not Detected/Quantitative Data R - Unusable Data
Q - Present in associated blanks (Semi-quantitative Data) UL - Not Detected/Biased Low Q - Not Applicable

AR302787

SUMMARY OF CLP ORGANIC ANALYSES - PHASE 1A
BERKS LANDFILL REMEDIAL INVESTIGATION
BERKS COUNTY, PENNSYLVANIA

Matrix: Groundwater Well

Semivolatile Organics

CROL	Parameter	Sample Point GM18S/1A Date Sampled: 09/03/92 Dilution Factor: Percent Moisture: NA %			Sample Point GM18S/1A Date Sampled: 09/03/92 Dilution Factor: Percent Moisture: NA %			Sample Point GM18S/RE Date Sampled: 09/03/92 Dilution Factor: Percent Moisture: NA %			Sample Point GUC30/1A Date Sampled: 09/04/92 Dilution Factor: Percent Moisture: NA %		
		SQL	Result	Qual	SQL	Result	Qual	SQL	Result	Qual	SQL	Result	Qual
10	Acenaphthene	10	10	U	10	10	U	10	10	U	10	10	U
10	Acenaphthylene	10	10	U	10	10	U	10	10	U	10	10	U
10	Anthracene	10	10	U	10	10	U	10	10	U	10	10	U
10	Benzo(a)Anthracene	10	10	U	10	10	U	10	10	U	10	10	U
10	4-Bromophenyl-phenylether	10	10	U	10	10	U	10	10	U	10	10	U
10	Butylbenzylphthalate	10	10	U	10	10	U	10	10	U	10	10	U
10	Carbazole	10	10	U	10	10	U	10	10	U	10	10	U
10	4-Chloro-3-Methylphenol	10	10	U	10	10	U	10	10	U	10	10	U
10	4-Chloroaniline	10	10	U	10	10	U	10	10	U	10	10	U
10	bis(2-Chloroethoxy)Methane	10	10	U	10	10	U	10	10	U	10	10	U
10	bis(2-Chloroethyl)ether	10	10	U	10	10	U	10	10	U	10	10	U
10	2-Chloronaphthalene	10	10	U	10	10	U	10	10	U	10	10	U
10	2-Chlorophenol	10	10	U	10	10	U	10	10	U	10	10	U
10	4-Chlorophenyl-phenylether	10	10	U	10	10	U	10	10	U	10	10	U
10	Chrysene	10	10	U	10	10	U	10	10	U	10	10	U
10	Dibenz(a,h)Anthracene	10	10	U	10	10	U	10	10	U	10	10	U
10	Dibenzofuran	10	10	U	10	10	U	10	10	U	10	10	U
10	1,2-Dichlorobenzene	10	10	U	10	10	U	10	10	U	10	10	U
10	1,3-Dichlorobenzene	10	10	U	10	10	U	10	10	U	10	10	U
10	1,4-Dichlorobenzene	10	10	U	10	10	U	10	10	U	10	10	U
10	3,3'-Dichlorobenzidine	10	10	U	10	10	U	10	10	U	10	10	U
10	2,4-Dichlorophenol	10	10	U	10	10	U	10	10	U	10	10	U
10	Diethylphthalate	10	10	U	10	10	U	10	10	U	10	10	U
10	2,4-Dimethylphenol	10	10	U	10	10	U	10	10	U	10	10	U
10	Dimethyl Phthalate	10	10	U	10	10	U	10	10	U	10	10	U
10	Di-n-butylphthalate	10	10	U	10	10	U	10	10	U	10	10	U
25	4,6-Dinitro-2-methylphenol	25	25	U	25	25	U	25	25	U	25	25	U
25	2,4-Dinitrophenol	25	25	U	25	25	U	25	25	U	25	25	U
10	2,4-Dinitrotoluene	10	10	U	10	10	U	10	10	U	10	10	U
10	2,6-Dinitrotoluene	10	10	U	10	10	U	10	10	U	10	10	U
10	Di-n-Octyl Phthalate	10	10	U	10	10	U	10	10	U	10	10	U
10	bis(2-Ethylhexyl)Phthalate	10	12	A	10	7	J	10	12	J	10	11	B
10	Fluorenone	10	10	U	10	10	U	10	10	U	10	10	U
10	Benzo(b)Fluoranthene	10	10	U	10	10	U	10	10	U	10	10	U
10	Benzo(k)Fluoranthene	10	10	U	10	10	U	10	10	U	10	10	U
10	Fluorene	10	10	U	10	10	U	10	10	U	10	10	U
10	Hexachlorobenzene	10	10	U	10	10	U	10	10	U	10	10	U
10	Hexachlorocyclopentadiene	10	10	U	10	10	U	10	10	U	10	10	U
10	Hexachloroethane	10	10	U	10	10	U	10	10	U	10	10	U

Notes:

All units are ug/l
Sample Quantitation Limit (SQL) is CROL multiplied by the dilution factor and adjusted for percent moisture as appropriate.
The qual column indicates the qualifier applied to the result following data validation (see below).

Qualifiers:.....
A - Acceptable (Quantitative) Data
J - Estimated (Semi-quantitative) Data
B - Present in associated blanks (Semi-quantitative Data)
K - Biased High (Semi-quantitative) Data
L - Biased Low (Semi-quantitative Data)
UL - Not Detected/Biased Low
UJ - Not Detected/Estimated Data
U - Not Detected/Quantitative Data
UL - Not Detected/Biased Low
N - Tentative Identification
R - Unusable Data
Q - Not Applicable

Matrix: Groundwater Well

Semi-volatile Organics

CROL	Parameter	Sample Point GJ18S/1A			Sample Point GJ18S/RE			Sample Point GJ18S/1A		
		SQL	Result	Qual	SQL	Result	Qual	SQL	Result	Qual
10	Indeno(1,2,3-cd)Pyrene	10	10	U	10	10	UJ	10	10	U
10	Isophorone	10	10	U	10	10	U	10	10	U
10	2-Methylnaphthalene	10	10	U	10	10	U	10	10	U
10	2-Methylphenol	10	10	U	10	10	U	10	10	U
10	4-Methylphenol	10	10	U	10	10	U	10	10	U
10	Naphthalene	10	10	U	10	10	U	10	10	U
25	2-Nitroaniline	25	25	U	25	25	U	25	25	UJ
25	3-Nitroaniline	25	25	U	25	25	U	25	25	U
25	4-Nitroaniline	25	25	U	25	25	U	25	25	U
10	Nitrobenzene	10	10	U	10	10	U	10	10	U
10	2-Nitrophenol	10	10	U	10	10	U	10	10	U
25	4-Nitrophenol	25	25	U	25	25	U	25	25	U
10	N-Nitroso-di-n-propylamine	10	10	U	10	10	U	10	10	U
10	N-Nitrosodiphenylamine	10	10	U	10	10	U	10	10	U
10	2,2'-Oxybis(1-Chloropropane)	10	10	U	10	10	U	10	10	U
25	Pentachlorophenol	25	25	U	25	25	U	25	25	U
10	Benzo(g,h,i)Perylene	10	10	UJ	10	10	UJ	10	10	U
10	Phenanthrene	10	10	U	10	10	U	10	10	U
10	Pyrene	10	10	U	10	10	U	10	10	U
10	Benzo(a)Pyrene	10	10	U	10	10	U	10	10	U
10	1,2,4-Trichlorobenzene	10	10	UJ	10	10	UJ	10	10	U
25	2,4,5-Trichlorophenol	25	25	U	25	25	U	25	25	U
10	2,4,6-Trichlorophenol	10	10	U	10	10	U	10	10	U

Notes:
All units are ug/l
Sample Quantitation Limit (SQL) is CROL multiplied by the dilution factor and adjusted for percent moisture as appropriate.
The dual column indicates the qualifier applied to the result following data validation (see below).

Qualifiers:.....
A - Acceptable (Quantitative) Data
J - Estimated (Semi-quantitative) Data
B - Present in associated blanks (Semi-quantitative Data)
K - Biased High (Semi-quantitative) Data
L - Biased Low (Semi-quantitative) Data
UJ - Not Detected/Estimated Data
U - Not Detected/Quantitative Data
UL - Not Detected/Biased Low
N - Tentative Identification
R - Unusable Data
Q - Not Applicable

SUMMARY OF CLP ORGANICS ANALYSES - PHASE 1A
BERKS LANDFILL REMEDIAL INVESTIGATION
BERKS COUNTY, PENNSYLVANIA

Matrix: Groundwater Well

Semivolatile Organics

CROL	Parameter	Sample Point GVC3S/1A			Sample Point GVC01/1A			Sample Point GVC02/1A			Sample Point GVC03/1A		
		SQL	Result	Qual	SQL	Result	Qual	SQL	Result	Qual	SQL	Result	Qual
10	Acenaphthene	10	10	U	10	10	R	10	10	U	10	10	U
10	Acenaphthylene	10	10	U	10	10	R	10	10	U	10	10	U
10	Anthracene	10	10	U	10	10	R	10	10	U	10	10	U
10	Benzo(a)Anthracene	10	10	U	10	10	R	10	10	U	10	10	U
10	4-Bromophenyl-phenylether	10	10	U	10	10	R	10	10	U	10	10	U
10	Butylbenzylphthalate	10	10	U	10	10	R	10	10	U	10	10	U
10	Carbazole	10	10	U	10	10	R	10	10	U	10	10	U
10	4-Chloro-3-Methylphenol	10	10	U	10	10	UL	10	10	U	10	10	U
10	4-Chloroaniline	10	10	U	10	10	R	10	10	U	10	10	U
10	bis(2-Chloroethoxy)Methane	10	10	U	10	10	R	10	10	U	10	10	U
10	bis(2-Chloroethyl)ether	10	10	U	10	10	R	10	10	U	10	10	U
10	2-Chloronaphthalene	10	10	U	10	10	R	10	10	U	10	10	U
10	2-Chlorophenol	10	10	U	10	10	UL	10	10	U	10	10	U
10	4-Chlorophenyl-phenylether	10	10	U	10	10	R	10	10	U	10	10	U
10	Chrysene	10	10	U	10	10	R	10	10	U	10	10	U
10	Dibenz(a,h)Anthracene	10	10	U	10	10	R	10	10	U	10	10	U
10	Dibenzofuran	10	10	U	10	10	R	10	10	U	10	10	U
10	1,2-Dichlorobenzene	10	10	U	10	10	R	10	10	U	10	10	U
10	1,3-Dichlorobenzene	10	10	U	10	10	R	10	10	U	10	10	U
10	1,4-Dichlorobenzene	10	10	U	10	10	R	10	10	U	10	10	U
10	3,3'-Dichlorobenzidine	10	10	U	10	10	R	10	10	U	10	10	U
10	2,4-Dichlorophenol	10	10	U	10	10	UL	10	10	U	10	10	U
10	Diethylphthalate	10	10	U	10	10	R	10	10	U	10	10	U
10	2,4-Dimethylphenol	10	10	U	10	10	UL	10	10	U	10	10	U
10	Dimethyl Phthalate	10	10	U	10	10	UL	10	10	U	10	10	U
10	Di-n-butylphthalate	10	10	U	10	10	R	10	10	U	10	10	U
25	4,6-Dinitro-2-methylphenol	25	25	U	25	25	UL	25	25	U	25	25	U
25	2,4-Dinitrophenol	25	25	U	25	25	UL	25	25	U	25	25	U
10	2,4-Dinitrotoluene	10	10	U	10	10	R	10	10	U	10	10	U
10	2,6-Dinitrotoluene	10	10	U	10	10	R	10	10	U	10	10	U
10	Di-n-Octyl Phthalate	10	10	UJ	10	10	R	10	10	U	10	10	U
10	bis(2-Ethylhexyl)Phthalate	10	5	J	10	10	R	10	10	U	10	1	B
10	Fluoranthene	10	10	U	10	10	R	10	10	U	10	10	U
10	Benzo(b)Fluoranthene	10	10	U	10	10	R	10	10	U	10	10	U
10	Benzo(k)Fluoranthene	10	10	U	10	10	R	10	10	U	10	10	U
10	Fluorene	10	10	U	10	10	R	10	10	U	10	10	U
10	Hexachlorobenzene	10	10	U	10	10	R	10	10	U	10	10	U
10	Hexachlorobutadiene	10	10	U	10	10	R	10	10	U	10	10	U
10	Hexachlorocyclopentadiene	10	10	U	10	10	R	10	10	U	10	10	U
10	Hexachloroethane	10	10	U	10	10	R	10	10	U	10	10	U

Notes:

All units are ug/l
Sample Quantitation Limit (SQL) is CROL multiplied by the dilution factor and adjusted for percent moisture as appropriate.
The Qual column indicates the qualifier applied to the result following data validation (see below).

Qualifiers:.....
A - Acceptable (Quantitative) Data
K - Biased High (Semi-quantitative) Data
J - Estimated (Semi-quantitative) Data
L - Biased Low (Semi-quantitative) Data
B - Present in associated blanks (Semi-quantitative) Data
UL - Not Detected/Estimated Data
UL - Not Detected/Quantitative Data
UL - Not Detected/Biased Low
N - Tentative Identification
R - Unusable Data
Q - Not Applicable

AR302790

Matrix: Groundwater Well

Semivolatile Organics

CRQL	Parameter	Sample Point GUC35/1A			Sample Point GUC01/1A			Sample Point GUC02/1A			Sample Point GUC03/1A		
		SQL	Result	Qual	SQL	Result	Qual	SQL	Result	Qual	SQL	Result	Qual
10	Indeno(1,2,3-cd)Pyrene	10	10	U	10	10	R	10	10	U	10	10	U
10	Isophorone	10	10	U	10	10	R	10	10	U	10	10	U
10	2-Methylnaphthalene	10	10	U	10	10	R	10	10	U	10	10	U
10	2-Methylphenol	10	10	U	10	10	UL	10	10	U	10	10	U
10	4-Methylphenol	10	10	U	10	10	UL	10	10	U	10	10	U
10	Naphthalene	10	10	U	10	10	R	10	10	U	10	10	U
25	2-Nitroaniline	25	25	U	25	25	R	25	25	U	25	25	U
25	3-Nitroaniline	25	25	U	25	25	R	25	25	U	25	25	U
25	4-Nitroaniline	25	25	U	25	25	R	25	25	U	25	25	U
10	Nitrobenzene	10	10	U	10	10	R	10	10	U	10	10	U
10	2-Nitrophenol	10	10	U	10	10	UL	10	10	U	10	10	U
25	4-Nitrophenol	25	25	U	25	25	UL	25	25	U	25	25	U
10	N-Nitroso-di-n-propylamine	10	10	U	10	10	R	10	10	U	10	10	U
10	N-Nitrosodiphenylamine	10	10	U	10	10	R	10	10	U	10	10	U
10	2,2'-Oxybis(1-Chloropropane)	10	10	U	10	10	R	10	10	U	10	10	U
25	Pentachlorophenol	25	25	U	25	25	UL	25	25	U	25	25	U
10	Benzof(g,h,i)perylene	10	10	U	10	10	R	10	10	U	10	10	U
10	Phenanthrene	10	10	U	10	10	R	10	10	U	10	10	U
10	Pyrene	10	10	U	10	10	UL	10	10	U	10	10	U
10	Benzof(a)Pyrene	10	10	U	10	10	R	10	10	U	10	10	U
10	1,2,6-Trichlorobenzene	10	10	U	10	10	R	10	10	U	10	10	U
25	2,4,5-Trichlorophenol	25	25	U	25	25	UL	25	25	U	25	25	U
10	2,4,6-Trichlorophenol	10	10	U	10	10	UL	10	10	U	10	10	U

Notes:
All units are ug/l
Sample Quantitation Limit (SQL) is CRQL multiplied by the dilution factor and adjusted for percent moisture as appropriate.
The Qual column indicates the qualifier applied to the result following data validation (see below).

Qualifiers:
A - Acceptable (Quantitative) Data
K - Biased High (Semi-quantitative) Data
J - Estimated (Semi-quantitative) Data
L - Biased Low (Semi-quantitative) Data
B - Present in associated blanks (Semi-quantitative Data)
U - Not Detected/Estimated Data
UL - Not Detected/Quantitative Data
R - Unusable Data
Q - Not Applicable
M - Tentative Identification

AR302791

Matrix: Groundwater Well

Semivolatile Organics

CRQL	Parameter	Sample Point GMC01/1B Date Sampled: 07/08/93 Dilution Factor: 1.0 Percent Moisture: NA %			Sample Point GMC02/1B Date Sampled: 07/08/93 Dilution Factor: 1.0 Percent Moisture: NA %			Sample Point GMC03B/1B Date Sampled: 07/06/93 Dilution Factor: 1.0 Percent Moisture: NA %			Sample Point GMC03B/FD Date Sampled: 07/06/93 Dilution Factor: 1.0 Percent Moisture: NA %		
		SOL	Result	Qual	SOL	Result	Qual	SOL	Result	Qual	SOL	Result	Qual
10	Acenaphthene	10	10	U	10	10	U	10	10	U	10	10	U
10	Acenaphthylene	10	10	U	10	10	U	10	10	U	10	10	U
10	Anthracene	10	10	U	10	10	U	10	10	U	10	10	U
10	Benzo(a)Anthracene	10	10	U	10	10	U	10	10	U	10	10	U
10	4-Bromophenyl-phthalate	10	10	U	10	10	U	10	10	U	10	10	U
10	Butylbenzylphthalate	10	10	U	10	10	U	10	10	U	10	10	U
10	Carbazole	10	10	U	10	10	U	10	10	U	10	10	U
10	4-Chloro-3-Methylphenol	10	10	U	10	10	U	10	10	U	10	10	U
10	4-Chloroaniline	10	10	U	10	10	U	10	10	U	10	10	U
10	bis(2-Chloroethoxy)Methane	10	10	U	10	10	U	10	10	U	10	10	U
10	bis(2-Chloroethyl)ether	10	10	U	10	10	U	10	10	U	10	10	U
10	2-Chloronaphthalene	10	10	U	10	10	U	10	10	U	10	10	U
10	2-Chlorophenol	10	10	U	10	10	U	10	10	U	10	10	U
10	4-Chlorophenyl-phenylether	10	10	U	10	10	U	10	10	U	10	10	U
10	Chrysene	10	10	U	10	10	U	10	10	U	10	10	U
10	Dibenz(a,h)Anthracene	10	10	U	10	10	U	10	10	U	10	10	U
10	Dibenzofuran	10	10	U	10	10	U	10	10	U	10	10	U
10	1,2-Dichlorobenzene	10	10	U	10	10	U	10	10	U	10	10	U
10	1,3-Dichlorobenzene	10	10	U	10	10	U	10	10	U	10	10	U
10	1,4-Dichlorobenzene	10	10	U	10	10	U	10	10	U	10	10	U
10	3,3'-Dichlorobenzidine	10	10	U	10	10	U	10	10	U	10	10	U
10	2,4-Dichlorophenol	10	10	U	10	10	U	10	10	U	10	10	U
10	Diethylphthalate	10	10	U	10	10	U	10	10	U	10	10	U
10	2,4-Dimethylphenol	10	10	U	10	10	U	10	10	U	10	10	U
10	Dimethyl-Phthalate	10	10	U	10	10	U	10	10	U	10	10	U
10	DJ-n-butylphthalate	10	10	U	10	10	U	10	10	U	10	10	U
25	4,6-Dinitro-2-methylphenol	25	25	U	25	25	U	25	25	U	25	25	U
25	2,4-Dinitrophenol	25	25	U	25	25	U	25	25	U	25	25	U
10	2,4-Dinitrotoluene	10	10	U	10	10	U	10	10	U	10	10	U
10	2,6-Dinitrotoluene	10	10	U	10	10	U	10	10	U	10	10	U
10	Di-n-Octyl Phthalate	10	10	U	10	10	U	10	10	U	10	10	U
10	bis(2-Ethylhexyl)Phthalate	10	10	U	10	10	U	10	10	U	10	10	U
10	Fluoranthene	10	10	U	10	10	U	10	10	U	10	10	U
10	Benzo(b)Fluoranthene	10	10	U	10	10	U	10	10	U	10	10	U
10	Benzo(k)Fluoranthene	10	10	U	10	10	U	10	10	U	10	10	U
10	Fluorene	10	10	U	10	10	U	10	10	U	10	10	U
10	Hexachlorobenzene	10	10	U	10	10	U	10	10	U	10	10	U
10	Hexachlorobutadiene	10	10	U	10	10	U	10	10	U	10	10	U
10	Hexachlorocyclopentadiene	10	10	U	10	10	U	10	10	U	10	10	U
10	Hexachloroethane	10	10	U	10	10	U	10	10	U	10	10	U

Notes:

All units are ug/l
Sample Quantitation Limit (SOL) is CRQL multiplied by the dilution factor and adjusted for percent moisture as appropriate.
The Qual column indicates the qualifier applied to the result following data validation (see below).

Qualifiers:.....
A - Acceptable (Quantitative) Data
J - Estimated (Semi-quantitative) Data
B - Present in associated blanks (Semi-quantitative Data)
K - Biased High (Semi-quantitative) Data
L - Biased Low (Semi-quantitative) Data
UJ - Not Detected/Estimated Data
U - Not Detected/Quantitative Data
UL - Not Detected/Biased Low
M - Tentative Identification
N - Unusable Data
R - Not Applicable
Q - Not Applicable

AR302792

Matrix: Groundwater Well

Semivolatile Organics

CRQL	Parameter	Sample Point GUC01/1B			Sample Point GUC02/1B			Sample Point GUC03D/1B			Sample Point GUC03D/FD		
		SQL	Result	Qual	SQL	Result	Qual	SQL	Result	Qual	SQL	Result	Qual
10	Indeno(1,2,3-cd)Pyrene	10	10	U	10	10	U	10	10	U	10	10	U
10	Isophorene	10	10	U	10	10	U	10	10	U	10	10	U
10	2-Methylnaphthalene	10	10	U	10	10	U	10	10	U	10	10	U
10	2-Methylphenol	10	10	U	10	10	U	10	10	U	10	10	U
10	4-Methylphenol	10	10	U	10	10	U	10	10	U	10	10	U
10	Naphthalene	10	10	U	10	10	U	10	10	U	10	10	U
25	2-Nitroaniline	25	25	U	25	25	U	25	25	U	25	25	U
25	3-Nitroaniline	25	25	U	25	25	U	25	25	U	25	25	U
25	4-Nitroaniline	25	25	U	25	25	U	25	25	U	25	25	U
10	Nitrobenzene	10	10	U	10	10	U	10	10	U	10	10	U
10	2-Nitrophenol	10	10	U	10	10	U	10	10	U	10	10	U
25	4-Nitrophenol	25	25	U	25	25	U	25	25	U	25	25	U
10	N-Nitroso-di-n-propylamine	10	10	U	10	10	U	10	10	U	10	10	U
10	N-Nitrosodiphenylamine	10	10	U	10	10	U	10	10	U	10	10	U
10	2,2'-Oxybis(1-Chloropropane)	10	10	U	10	10	U	10	10	U	10	10	U
25	Pentachlorophenol	25	25	U	25	25	U	25	25	U	25	25	U
10	Benzo(g,h,i)perylene	10	10	U	10	10	U	10	10	U	10	10	U
10	Phenanthrene	10	10	U	10	10	U	10	10	U	10	10	U
10	Phenol	10	10	U	10	10	U	10	10	U	10	10	U
10	Pyrene	10	10	U	10	10	U	10	10	U	10	10	U
10	Benzo(a)Pyrene	10	10	U	10	10	U	10	10	U	10	10	U
10	1,2,4-Trichlorobenzene	10	10	U	10	10	U	10	10	U	10	10	U
25	2,4,5-Trichlorophenol	25	25	U	25	25	U	25	25	U	25	25	U
10	2,4,6-Trichlorophenol	10	10	U	10	10	U	10	10	U	10	10	U

Notes:

All units are ug/l
Sample Quantitation Limit (SQL) is CRQL multiplied by the dilution factor and adjusted for percent moisture as appropriate.
The Qual column indicates the qualifier applied to the result following data validation (see below).

Qualifiers:.....
A - Acceptable (Quantitative) Data K - Biased High (Semi-quantitative) Data UJ - Not Detected/Estimated Data N - Tentative Identification
J - Estimated (Semi-quantitative) Data L - Biased Low (Semi-quantitative) Data U - Not Detected/Quantitative Data R - Unusable Data
UL - Present in associated blanks (Semi-quantitative Data) UL - Not Detected/Biased Low UL - Not Detected/Biased Low Q - Not Applicable

AR302793

SUMMARY OF CLP ORGANIC ANALYSES - PHASE 1B
BERKS LANDFILL REMEDIAL INVESTIGATION
BERKS COUNTY, PENNSYLVANIA

Matrix: Groundwater Well

Semivolatiles Organics

CRQL	Parameter	Sample Point GVC03D/FR Date Sampled: 07/06/93 Dilution Factor: 1.0 Percent Moisture: NA %			Sample Point GVC03S/1B Date Sampled: 07/02/93 Dilution Factor: 1.0 Percent Moisture: NA %			Sample Point GVC04D/1B Date Sampled: 06/30/93 Dilution Factor: 1.0 Percent Moisture: NA %			Sample Point GVC04S/1B Date Sampled: 06/30/93 Dilution Factor: 1.0 Percent Moisture: NA %		
		SQL	Result	Qual	SQL	Result	Qual	SQL	Result	Qual	SQL	Result	Qual
10	Acenaphthene	10	10	U	10	10	U	10	10	U	10	10	U
10	Acenaphthylene	10	10	U	10	10	U	10	10	U	10	10	U
10	Anthracene	10	10	U	10	10	U	10	10	U	10	10	U
10	Benzo(a)Anthracene	10	10	UJ	10	10	U	10	10	U	10	10	U
10	4-Bromophenyl-phenylether	10	10	U	10	10	U	10	10	U	10	10	U
10	Butylbenzylphthalate	10	10	UJ	10	10	U	10	10	U	10	10	U
10	Carbazole	10	10	U	10	10	U	10	10	U	10	10	U
10	4-Chloro-3-Methylphenol	10	10	U	10	10	U	10	10	U	10	10	U
10	4-Chloroaniline	10	10	U	10	10	U	10	10	U	10	10	U
10	bis(2-Chloroethoxy)Methane	10	10	U	10	10	U	10	10	U	10	10	U
10	bis(2-Chloroethyl)ether	10	10	U	10	10	U	10	10	U	10	10	U
10	2-Chloronaphthalene	10	10	U	10	10	U	10	10	U	10	10	U
10	2-Chlorophenol	10	10	U	10	10	U	10	10	U	10	10	U
10	4-Chlorophenyl-phenylether	10	10	U	10	10	U	10	10	U	10	10	U
10	Chrysene	10	10	UJ	10	10	U	10	10	U	10	10	U
10	DiBenz(a,h)Anthracene	10	10	UJ	10	10	U	10	10	U	10	10	U
10	Dibenzofuran	10	10	U	10	10	U	10	10	U	10	10	U
10	1,2-Dichlorobenzene	10	10	U	10	10	U	10	10	U	10	10	U
10	1,3-Dichlorobenzene	10	10	U	10	10	U	10	10	U	10	10	U
10	1,4-Dichlorobenzene	10	10	U	10	10	U	10	10	U	10	10	U
10	3,3'-Dichlorobenzidine	10	10	U	10	10	U	10	10	U	10	10	U
10	2,4-Dichlorophenol	10	10	UJ	10	10	U	10	10	U	10	10	U
10	Diethylphthalate	10	10	U	10	10	U	10	10	U	10	10	U
10	2,4-Dimethylphenol	10	10	B	10	10	U	10	10	U	10	10	U
10	Dimethyl Phthalate	10	10	U	10	10	U	10	10	U	10	10	U
10	Di-n-butylphthalate	10	10	B	10	10	U	10	10	U	10	10	U
25	4,6-Dinitro-2-methylphenol	25	25	U	25	25	U	25	25	U	25	25	U
25	2,4-Dinitrophenol	25	25	U	25	25	U	25	25	U	25	25	U
10	2,4-Dinitrotoluene	10	10	U	10	10	U	10	10	U	10	10	U
10	2,6-Dinitrotoluene	10	10	U	10	10	U	10	10	U	10	10	U
10	Di-n-Octyl Phthalate	10	10	UJ	10	10	U	10	10	U	10	10	U
10	bis(2-Ethylhexyl)Phthalate	10	10	B	10	10	U	10	10	U	10	10	U
10	Fluoranthene	10	10	U	10	10	U	10	10	U	10	10	U
10	Benzo(b)Fluoranthene	10	10	UJ	10	10	U	10	10	U	10	10	U
10	Benzo(k)Fluoranthene	10	10	UJ	10	10	U	10	10	U	10	10	U
10	Fluorene	10	10	U	10	10	U	10	10	U	10	10	U
10	Hexchlorobenzene	10	10	U	10	10	U	10	10	U	10	10	U
10	Hexchlorobutadiene	10	10	U	10	10	U	10	10	U	10	10	U
10	Hexchlorocyclopentadiene	10	10	U	10	10	U	10	10	U	10	10	U
10	Hexchloroethene	10	10	U	10	10	U	10	10	U	10	10	U

Notes:

All units are ug/l
Sample Quantitation Limit (SQL) is CRQL multiplied by the dilution factor and adjusted for percent moisture as appropriate.
The Qual column indicates the qualifier applied to the result following data validation (see below).

Qualifiers:
A - Acceptable (Quantitative) Data
J - Estimated (Semi-quantitative) Data
B - Present in associated blanks (Semi-quantitative Data)
K - Biased High (Semi-quantitative) Data
L - Biased Low (Semi-quantitative) Data
UJ - Not Detected/Estimated Data
U - Not Detected/Quantitative Data
UL - Not Detected/Biased Low
N - Tentative Identification
R - Unusable Data
Q - Not Applicable

AR302794

Matrix: Groundwater Well

Semi-volatile Organics

CROL	Parameter	Sample Point GUC030/FR Date Sampled: 07/06/93 Dilution Factor: 1.0 Percent Moisture: NA %			Sample Point GUC03S/19 Date Sampled: 07/02/93 Dilution Factor: 1.0 Percent Moisture: NA %			Sample Point GUC040/18 Date Sampled: 06/30/93 Dilution Factor: 1.0 Percent Moisture: NA %			Sample Point GUC04S/18 Date Sampled: 06/30/93 Dilution Factor: 1.0 Percent Moisture: NA %		
		SQL	Result	Qual	SQL	Result	Qual	SQL	Result	Qual	SQL	Result	Qual
10	Indeno(1,2,3-cd)Pyrene	10	10	UJ	10	10	U	10	10	10	10	U	
10	Isophorone	10	10	U	10	10	U	10	10	10	10	U	
10	2-Methylnaphthalene	10	10	U	10	10	U	10	10	10	10	U	
10	2-Methylphenol	10	10	U	10	10	U	10	10	10	10	U	
10	4-Methylphenol	10	10	U	10	10	U	10	10	10	10	U	
10	Naphthalene	10	10	U	10	10	U	10	10	10	10	U	
25	2-Nitroaniline	25	25	U	25	25	U	25	25	25	25	U	
25	3-Nitroaniline	25	25	U	25	25	U	25	25	25	25	U	
25	4-Nitroaniline	25	25	U	25	25	U	25	25	25	25	U	
10	Nitrobenzene	10	10	U	10	10	U	10	10	10	10	U	
10	2-Nitrophenol	10	10	U	10	10	U	10	10	10	10	U	
25	4-Nitrophenol	25	25	U	25	25	U	25	25	25	25	U	
10	N-Nitroso-di-n-propylamine	10	10	U	10	10	U	10	10	10	10	U	
10	N-Nitrosodiphenylamine	10	10	U	10	10	U	10	10	10	10	U	
10	2,2'-Oxybis(1-Chloropropane)	10	10	U	10	10	U	10	10	10	10	U	
25	Pentachlorophenol	25	25	UJ	25	25	U	25	25	25	25	U	
10	Benzof(g,h,i)Perylene	10	10	U	10	10	U	10	10	10	10	U	
10	Phenanthrene	10	10	U	10	10	U	10	10	10	10	U	
10	Phenol	10	4	B	10	10	U	10	3	10	10	U	
10	Pyrene	10	10	UJ	10	10	U	10	10	10	10	U	
10	Benzof(a)Pyrene	10	10	UJ	10	10	U	10	10	10	10	U	
10	1,2,6-Trichlorobenzene	10	10	U	10	10	U	10	10	10	10	U	
25	2,4,5-Trichlorophenol	25	25	U	25	25	U	25	25	25	25	U	
10	2,4,6-Trichlorophenol	10	10	U	10	10	U	10	10	10	10	U	

Notes:
All units are ug/l
Sample Quantitation Limit (SQL) is CROL multiplied by the dilution factor and adjusted for percent moisture as appropriate.
The Qual column indicates the qualifier applied to the result following data validation (see below).

Qualifiers:
A - Acceptable (Quantitative) Data
K - Biased High (Semi-quantitative) Data
J - Estimated (Semi-quantitative) Data
L - Biased Low (Semi-quantitative) Data
% - Present in associated blanks (Semi-quantitative Data)
UJ - Not Detected/Estimated Data
U - Not Detected/Quantitative Data
UL - Not Detected/Biased Low
M - Tentative Identification
R - Unusable Data
B - Not Applicable

AR302795

SUMMARY OF CLP ORGANIC ANALYSES - PHASE 1B
BERKS LANDFILL REMEDIAL INVESTIGATION
BERKS COUNTY, PENNSYLVANIA

Matrix: Groundwater Well

Semivolatile Organics

CROL	Parameter	Sample Point GWC04S/RE Date Sampled: 06/30/93 Dilution Factor: 1.0 Percent Moisture: NA %			Sample Point GWC06D/1B Date Sampled: 07/12/93 Dilution Factor: 1.0 Percent Moisture: NA %			Sample Point GWC06S/1B Date Sampled: 07/12/93 Dilution Factor: 1.0 Percent Moisture: NA %			Sample Point GWC07D/1B Date Sampled: 07/09/93 Dilution Factor: 1.0 Percent Moisture: NA %		
		SQL	Result	Qual	SQL	Result	Qual	SQL	Result	Qual	SQL	Result	Qual
10	Acenaphthene	10	10	U	10	10	U	10	10	U	10	10	U
10	Acenaphthylene	10	10	U	10	10	U	10	10	U	10	10	U
10	Anthracene	10	10	U	10	10	U	10	10	U	10	10	U
10	Benzo(a)Anthracene	10	10	U	10	10	U	10	10	U	10	10	U
10	4-Bromophenyl-phenylether	10	10	U	10	10	U	10	10	U	10	10	U
10	Butylbenzylphthalate	10	10	U	10	10	U	10	10	U	10	10	U
10	Carbazole	10	10	U	10	10	U	10	10	U	10	10	U
10	4-Chloro-3-Methylphenol	10	10	R	10	10	U	10	10	U	10	10	U
10	4-Chloroaniline	10	10	U	10	10	U	10	10	U	10	10	U
10	bis(2-Chloroethoxy)Methane	10	10	U	10	10	U	10	10	U	10	10	U
10	bis(2-Chloroethyl)Ether	10	10	U	10	10	U	10	10	U	10	10	U
10	2-Chloronaphthalene	10	10	U	10	10	U	10	10	U	10	10	U
10	2-Chlorophenol	10	10	R	10	10	U	10	10	U	10	10	U
10	4-Chlorophenyl-phenylether	10	10	U	10	10	U	10	10	U	10	10	U
10	Chrysene	10	10	U	10	10	U	10	10	U	10	10	U
10	Dibenz(a,h)Anthracene	10	10	U	10	10	U	10	10	U	10	10	U
10	Dibenzofuran	10	10	U	10	10	U	10	10	U	10	10	U
10	1,2-Dichlorobenzene	10	10	U	10	10	U	10	10	U	10	10	U
10	1,3-Dichlorobenzene	10	10	U	10	10	U	10	10	U	10	10	U
10	1,4-Dichlorobenzene	10	10	U	10	10	U	10	10	U	10	10	U
10	3,3'-Dichlorobenzidine	10	10	U	10	10	U	10	10	U	10	10	U
10	2,4-Dichlorophenol	10	10	R	10	10	U	10	10	U	10	10	U
10	Diethylphthalate	10	10	U	10	10	U	10	10	U	10	10	U
10	2,4-Dimethylphenol	10	10	R	10	10	U	10	10	U	10	10	U
10	Dimethyl Phthalate	10	10	U	10	10	U	10	10	U	10	10	U
10	Di-n-butylphthalate	10	10	U	10	10	U	10	10	U	10	10	U
25	4,6-Dinitro-2-methylphenol	25	25	R	25	25	U	25	25	U	25	25	U
25	2,4-Dinitrophenol	25	25	R	25	25	U	25	25	U	25	25	U
10	2,4-Dinitrotoluene	10	10	U	10	10	U	10	10	U	10	10	U
10	2,6-Dinitrotoluene	10	10	U	10	10	U	10	10	U	10	10	U
10	Di-n-Octyl Phthalate	10	10	U	10	10	U	10	10	U	10	10	U
10	bis(2-Ethylhexyl)Phthalate	10	10	U	10	10	U	10	10	U	10	10	U
10	Fluoranthene	10	10	U	10	10	U	10	10	U	10	10	U
10	Benzo(b)Fluoranthene	10	10	U	10	10	U	10	10	U	10	10	U
10	Benzo(k)Fluoranthene	10	10	U	10	10	U	10	10	U	10	10	U
10	Fluorene	10	10	U	10	10	U	10	10	U	10	10	U
10	Hexachlorobenzene	10	10	U	10	10	U	10	10	U	10	10	U
10	Hexachlorobutadiene	10	10	U	10	10	U	10	10	U	10	10	U
10	Hexachlorocyclopentadiene	10	10	U	10	10	U	10	10	U	10	10	U
10	Hexachloroethane	10	10	U	10	10	U	10	10	U	10	10	U

Notes:

All units are ug/l
Sample Quantitation Limit (SQL) is CROL multiplied by the dilution factor and adjusted for percent moisture as appropriate.
The Qual column indicates the qualifier applied to the result following data validation (see below).

Qualifiers:
A - Acceptable (Quantitative) Data K - Biased High (Semi-quantitative) Data UJ - Not Detected/Estimated Data N - Tentative Identification
J - Estimated (Semi-quantitative) Data L - Biased Low (Semi-quantitative) Data U - Not Detected/Quantitative Data R - Unusable Data
B - Present in associated blanks (Semi-quantitative Data) UL - Not Detected/Biased Low Q - Not Applicable

AR302796

Matrix: Groundwater Well

semi-volatile Organics

CRQL	Parameter	Sample Point GUC04S/BE			Sample Point GUC06D/1B			Sample Point GUC06S/1B			Sample Point GUC07D/1B		
		SQL	Result	Qual	SQL	Result	Qual	SQL	Result	Qual	SQL	Result	Qual
10	Indeno(1,2,3-cd)Pyrene	10	10	U	10	10	U	10	10	U	10	10	U
10	Isophorone	10	10	U	10	10	U	10	10	U	10	10	U
10	2-Methylnaphthalene	10	10	U	10	10	U	10	10	U	10	10	U
10	2-Methylphenol	10	10	R	10	10	U	10	10	U	10	10	U
10	4-Methylphenol	10	10	R	10	10	U	10	10	U	10	10	U
10	Naphthalene	10	10	U	10	10	U	10	10	U	10	10	U
25	2-Nitroaniline	25	25	U	25	25	U	25	25	U	25	25	U
25	3-Nitroaniline	25	25	U	25	25	U	25	25	U	25	25	U
25	4-Nitroaniline	25	25	U	25	25	U	25	25	U	25	25	U
10	Nitrobenzene	10	10	U	10	10	U	10	10	U	10	10	U
10	2-Nitrophenol	10	10	R	10	10	U	10	10	U	10	10	U
25	4-Nitrophenol	25	25	R	25	25	U	25	25	U	25	25	U
10	N-Nitroso-di-n-propylamine	10	10	U	10	10	U	10	10	U	10	10	U
10	N-Nitrosodiphenylamine	10	10	U	10	10	U	10	10	U	10	10	U
10	2,2'-Oxybis(1-Chloropropane)	10	10	U	10	10	U	10	10	U	10	10	U
25	Pentachloroethene	25	25	U	25	25	U	25	25	U	25	25	U
10	Benzo(g,h,i)Perylene	10	10	U	10	10	U	10	10	U	10	10	U
10	Phenanthrene	10	10	U	10	10	U	10	10	U	10	10	U
10	Pyrene	10	10	U	10	10	U	10	10	U	10	10	U
10	Benzo(a)Pyrene	10	10	U	10	10	U	10	10	U	10	10	U
10	1,2,4-Trichlorobenzene	10	10	U	10	10	U	10	10	U	10	10	U
25	2,4,5-Trichlorophenol	25	25	R	25	25	U	25	25	U	25	25	U
10	2,4,6-Trichlorophenol	10	10	R	10	10	U	10	10	U	10	10	U

Notes:
All units are ug/l
Sample Quantitation Limit (SQL) is CRQL multiplied by the dilution factor and adjusted for percent moisture as appropriate.
The qual column indicates the qualifier applied to the result following data validation (see below).

Qualifiers:
A - Acceptable (Quantitative) Data
K - Biased High (Semi-quantitative) Data
L - Biased Low (Semi-quantitative) Data
B - Present in associated blanks (Semi-quantitative Data)
J - Estimated (Semi-quantitative) Data
U - Not Detected/Quantitative Data
UL - Not Detected/Biased Low
M - Tentative Identification
R - Unusable Data
Q - Not Applicable

AR302797

Matrix: Groundwater Well

Semivolatile Organics

CROL	Parameter	Sample Point GMC07S/1B			Sample Point GMC02/1B			Sample Point GMC01/1B			Sample Point GMC02/1B		
		SQL	Result	Qual	SQL	Result	Qual	SQL	Result	Qual	SQL	Result	Qual
10	Acenaphthene	10	10	U	10	10	U	10	10	U	10	10	U
10	Acenaphthylene	10	10	U	10	10	U	10	10	U	10	10	U
10	Anthracene	10	10	U	10	10	U	10	10	U	10	10	U
10	Benzo(a)Anthracene	10	10	U	10	10	U	10	10	U	10	10	U
10	4-Bromophenyl-phthalate	10	10	U	10	10	U	10	10	U	10	10	U
10	Butylbenzylphthalate	10	2	J	10	10	U	10	10	U	10	10	U
10	Carbazole	10	10	U	10	10	U	10	10	U	10	10	U
10	4-Chloro-3-Methylphenol	10	10	U	10	10	U	10	10	U	10	10	U
10	4-Chloroaniline	10	10	U	10	10	U	10	10	U	10	10	U
10	bis(2-Chloroethoxy)Methane	10	10	U	10	10	U	10	10	U	10	10	U
10	bis(2-Chloroethyl)ether	10	10	U	10	10	U	10	10	U	10	10	U
10	2-Chloronaphthalene	10	10	U	10	10	U	10	10	U	10	10	U
10	2-Chlorophenol	10	10	U	10	10	U	10	10	U	10	10	U
10	4-Chlorophenyl-phenylether	10	10	U	10	10	U	10	10	U	10	10	U
10	Chrysene	10	10	U	10	10	U	10	10	U	10	10	U
10	Di-benz(a,h)Anthracene	10	10	U	10	10	U	10	10	U	10	10	U
10	Dibenzofuran	10	10	U	10	10	U	10	10	U	10	10	U
10	1,2-Dichlorobenzene	10	10	U	10	10	U	10	10	U	10	10	U
10	1,3-Dichlorobenzene	10	10	U	10	10	U	10	10	U	10	10	U
10	1,4-Dichlorobenzene	10	5	J	10	10	U	10	10	U	10	10	U
10	3,3'-Dichlorobenzidine	10	10	U	10	10	U	10	10	U	10	10	U
10	2,4-Dichlorophenol	10	10	U	10	10	U	10	10	U	10	10	U
10	Diethylphthalate	10	3	B	10	10	U	10	10	U	10	10	U
10	Dimethyl Phthalate	10	10	U	10	10	U	10	10	U	10	10	U
10	Di-n-butylphthalate	10	1	B	10	10	U	10	10	U	10	10	U
25	4,6-Dinitro-2-methylphenol	25	25	U	25	25	U	25	25	U	25	25	U
25	2,4-Dinitrophenol	25	25	U	25	25	U	25	25	U	25	25	U
10	2,4-Dinitrotoluene	10	10	U	10	10	U	10	10	U	10	10	U
10	2,6-Dinitrotoluene	10	10	U	10	10	U	10	10	U	10	10	U
10	Di-n-Octyl Phthalate	10	10	U	10	10	U	10	10	U	10	10	U
10	Bis(2-Ethylhexyl)Phthalate	10	4	B	10	10	U	10	10	U	10	10	U
10	Fluoranthene	10	10	U	10	10	U	10	10	U	10	10	U
10	Benzo(b)Fluoranthene	10	10	U	10	10	U	10	10	U	10	10	U
10	Benzo(k)Fluoranthene	10	10	U	10	10	U	10	10	U	10	10	U
10	Fluorene	10	10	U	10	10	U	10	10	U	10	10	U
10	Hexachlorobenzene	10	10	U	10	10	U	10	10	U	10	10	U
10	Hexachlorobutadiene	10	10	U	10	10	U	10	10	U	10	10	U
10	Hexachlorocyclopentadiene	10	10	U	10	10	U	10	10	U	10	10	U
10	Hexachloroethane	10	10	U	10	10	U	10	10	U	10	10	U

Notes:

All units are ug/l
Sample Quantitation Limit (SQL) is CROL multiplied by the dilution factor and adjusted for percent moisture as appropriate.
The Qual column indicates the qualifier applied to the result following data validation (see below).

Qualifiers:.....
A - Acceptable (Quantitative) Data
K - Biased High (Semi-quantitative) Data
UJ - Not Detected/Estimated Data
N - Tentative Identification
J - Estimated (Semi-quantitative) Data
L - Biased Low (Semi-quantitative) Data
UL - Not Detected/Quantitative Data
R - Unusable Data
B - Present in associated blanks (Semi-quantitative Data)
UL - Not Detected/Biased Low
Q - Not Applicable

AR302798

Matrix: Groundwater Well

Semivolatile Organics

CROL	Parameter	Sample Point GUC07S/18		Sample Point GUC02/18		Sample Point GUC01/18		Sample Point GUC02/18	
		SQL	Result	SQL	Result	SQL	Result	SQL	Result
10	Indeno(1,2,3-cd)Pyrene	10	10	10	10	10	10	10	10
10	Isophorone	10	10	10	10	10	10	10	10
10	2-Methylnaphthalene	10	10	10	10	10	10	10	10
10	2-Methylphenol	10	10	10	10	10	10	10	10
10	4-Methylphenol	10	10	10	10	10	10	10	10
25	Naphthalene	25	25	25	25	25	25	25	25
25	2-Nitroaniline	25	25	25	25	25	25	25	25
25	3-Nitroaniline	25	25	25	25	25	25	25	25
10	Nitrobenzene	10	10	10	10	10	10	10	10
10	2-Nitrophenol	10	10	10	10	10	10	10	10
25	4-Nitrophenol	25	25	25	25	25	25	25	25
10	N-Nitroso-di-n-propylamine	10	10	10	10	10	10	10	10
10	N-Nitrosodiphenylamine	10	10	10	10	10	10	10	10
10	2,2'-Oxybis(1-Chloropropane)	10	10	10	10	10	10	10	10
25	Pentachlorophenol	25	25	25	25	25	25	25	25
10	Benzo(g,h,i)perylene	10	10	10	10	10	10	10	10
10	Phenanthrene	10	10	10	10	10	10	10	10
10	Pyrene	10	3	10	10	10	10	10	10
10	Benzo(a)Pyrene	10	10	10	10	10	10	10	10
10	1,2,4-Trichlorobenzene	10	10	10	10	10	10	10	10
25	2,4,5-Trichlorophenol	25	25	25	25	25	25	25	25
10	2,4,6-Trichlorophenol	10	10	10	10	10	10	10	10

Notes:
All units are ug/l
Sample Quantitation Limit (SQL) is CROL multiplied by the dilution factor and adjusted for percent moisture as appropriate.
The Qual column indicates the qualifier applied to the result following data validation (see below).

Qualifiers:
A - Acceptable (Quantitative) Data
K - Biased High (Semi-quantitative) Data
L - Biased Low (Semi-quantitative) Data
J - Estimated (Semi-quantitative) Data
g - Present in associated blanks (Semi-quantitative Data)
UJ - Not Detected/Estimated Data
U - Not Detected/Quantitative Data
UL - Not Detected/Biased Low
M - Tentative Identification
N - Unusable Data
R - Not Applicable
Q - Not Applicable

AR302799

SUMMARY OF CLP ORGANICS ANALYSES - PHASE 1B
BERKS LANDFILL REMEDIAL INVESTIGATION
BERKS COUNTY, PENNSYLVANIA

Matrix: Groundwater Well

Semivolatile Organics

CROL	Parameter	Sample Point GM03/1B			Sample Point GM04/1B			Sample Point GM05/1B			Sample Point GM06/1B		
		SQL	Result	Qual	SQL	Result	Qual	SQL	Result	Qual	SQL	Result	Qual
10	Acenaphthene	10	10	U	10	10	U	10	10	U	10	10	U
10	Acenaphthylene	10	10	U	10	10	U	10	10	U	10	10	U
10	Anthracene	10	10	U	10	10	U	10	10	U	10	10	U
10	Benzo(a)Anthracene	10	10	U	10	10	U	10	10	U	10	10	U
10	4-Bromophenyl-phenylether	10	10	U	10	10	U	10	10	U	10	10	U
10	Butylbenzylphthalate	10	10	U	10	10	U	10	10	U	10	10	U
10	Carbazole	10	10	U	10	10	U	10	10	U	10	10	U
10	4-Chloro-3-Methylphenol	10	10	U	10	10	U	10	10	U	10	10	U
10	4-Chloroaniline	10	10	U	10	10	U	10	10	U	10	10	U
10	bis(2-Chloroethoxy)Methane	10	10	U	10	10	U	10	10	U	10	10	U
10	bis(2-Chloroethyl)ether	10	10	U	10	10	U	10	10	U	10	10	U
10	2-Chloronaphthalene	10	10	U	10	10	U	10	10	U	10	10	U
10	2-Chlorophenol	10	10	U	10	10	U	10	10	U	10	10	U
10	4-Chlorophenyl-phenylether	10	10	U	10	10	U	10	10	U	10	10	U
10	Chrysene	10	10	U	10	10	U	10	10	U	10	10	U
10	Dibenz(a,h)Anthracene	10	10	U	10	10	U	10	10	U	10	10	U
10	Dibenzofuran	10	10	U	10	10	U	10	10	U	10	10	U
10	1,2-Dichlorobenzene	10	10	U	10	10	U	10	10	U	10	10	U
10	1,3-Dichlorobenzene	10	10	U	10	10	U	10	10	U	10	10	U
10	1,4-Dichlorobenzene	10	10	U	10	10	U	10	10	U	10	10	U
10	3,3'-Dichlorobenzidine	10	10	U	10	10	U	10	10	U	10	10	U
10	2,4-Dichlorophenol	10	10	U	10	10	U	10	10	U	10	10	U
10	Diethylphthalate	10	10	U	10	10	U	10	10	U	10	10	U
10	2,4-Dimethylphenol	10	10	U	10	10	U	10	10	U	10	10	U
10	Dimethyl phthalate	10	10	U	10	10	U	10	10	U	10	10	U
10	Di-n-butylphthalate	10	10	U	10	10	U	10	10	U	10	10	U
25	4,6-Dinitro-2-methylphenol	25	25	U	25	25	U	25	25	U	25	25	U
25	2,4-Dinitrophenol	25	25	U	25	25	U	25	25	U	25	25	U
10	2,4-Dinitrotoluene	10	10	U	10	10	U	10	10	U	10	10	U
10	2,6-Dinitrotoluene	10	10	U	10	10	U	10	10	U	10	10	U
10	Di-n-Octyl Phthalate	10	10	U	10	10	U	10	10	U	10	10	U
10	bis(2-Ethylhexyl)Phthalate	10	2	U	10	18	U	10	5	U	10	10	U
10	Fluoranthene	10	10	U	10	10	U	10	10	U	10	10	U
10	Benzo(b)Fluoranthene	10	10	U	10	10	U	10	10	U	10	10	U
10	Benzo(k)Fluoranthene	10	10	U	10	10	U	10	10	U	10	10	U
10	Fluorene	10	10	U	10	10	U	10	10	U	10	10	U
10	Hexachlorobenzene	10	10	U	10	10	U	10	10	U	10	10	U
10	Hexachlorobutadiene	10	10	U	10	10	U	10	10	U	10	10	U
10	Hexachlorocyclopentadiene	10	10	U	10	10	U	10	10	U	10	10	U
10	Hexachloroethane	10	10	U	10	10	U	10	10	U	10	10	U

Notes:

All units are ug/l
Sample Quantitation Limit (SQL) is CROL multiplied by the dilution factor and adjusted for percent moisture as appropriate.
The Qual column indicates the qualifier applied to the result following data validation (see below).

Qualifiers:.....
A - Acceptable (Quantitative) Data
J - Estimated (Semi-quantitative) Data
B - Present in associated blanks (Semi-quantitative Data)
K - Biased High (Semi-quantitative) Data
L - Biased Low (Semi-quantitative) Data
UJ - Not Detected/Estimated Data
U - Not Detected/Quantitative Data
UL - Not Detected/Biased Low
M - Tentative Identification
N - Unusable Data
R - Unusable Data
Q - Not Applicable

AR302800

Matrix: Groundwater Well

Semivolatile Organics

CRQL	Parameter	Sample Point GUC03/18			Sample Point GUC04/18			Sample Point GUC05/18			Sample Point GUC06/18		
		SQL	Result	Qual	SQL	Result	Qual	SQL	Result	Qual	SQL	Result	Qual
10	Indeno(1,2,3-cd)Pyrene	10	10	U	10	10	U	10	10	U	10	10	U
10	Isophorone	10	10	U	10	10	U	10	10	U	10	10	U
10	2-Methylnaphthalene	10	10	U	10	10	U	10	10	U	10	10	U
10	2-Methylphenol	10	10	U	10	10	U	10	10	U	10	10	U
10	4-Methylphenol	10	10	U	10	10	U	10	10	U	10	10	U
10	Naphthalene	10	10	U	10	10	U	10	10	U	10	10	U
25	2-Nitroaniline	25	25	U	25	25	U	25	25	U	25	25	U
25	3-Nitroaniline	25	25	U	25	25	U	25	25	U	25	25	U
25	4-Nitroaniline	25	25	U	25	25	U	25	25	U	25	25	U
10	Nitrobenzene	10	10	U	10	10	U	10	10	U	10	10	U
10	2-Nitrophenol	10	10	U	10	10	U	10	10	U	10	10	U
25	4-Nitrophenol	25	25	U	25	25	U	25	25	U	25	25	U
10	N-Nitroso-di-n-propylamine	10	10	U	10	10	U	10	10	U	10	10	U
10	N-Nitrosodiphenylamine	10	10	U	10	10	U	10	10	U	10	10	U
10	2,2'-Oxybis(1-Chloropropane)	10	10	U	10	10	U	10	10	U	10	10	U
25	Pentachlorophenol	25	25	U	25	25	U	25	25	U	25	25	U
10	Benzo(g,h,i)Perylene	10	10	U	10	10	U	10	10	U	10	10	U
10	Phenanthrene	10	10	U	10	10	U	10	10	U	10	10	U
10	Phenol	10	10	U	10	10	U	10	10	U	10	10	U
10	Pyrene	10	10	U	10	10	U	10	10	U	10	10	U
10	Benzo(a)Pyrene	10	10	U	10	10	U	10	10	U	10	10	U
10	1,2,4-Trichlorobenzene	10	10	U	10	10	U	10	10	U	10	10	U
25	2,4,5-Trichlorophenol	25	25	U	25	25	U	25	25	U	25	25	U
10	2,4,6-Trichlorophenol	10	10	U	10	10	U	10	10	U	10	10	U

Notes:
All units are ug/l
Sample quantitation limit (SQL) is CRQL multiplied by the dilution factor and adjusted for percent moisture as appropriate.
The qual column indicates the qualifier applied to the result following data validation (see below).

Qualifiers:
A - Acceptable (Quantitative) Data K - Biased High (Semi-quantitative) Data JJ - Not Detected/Estimated Data M - Tentative Identification
J - Estimated (Semi-quantitative) Data L - Biased Low (Semi-quantitative) Data U - Not Detected/Quantitative Data R - Unusable Data
- Present in associated blanks (Semi-quantitative Data) UL - Not Detected/Biased Low Q - Not Applicable

AR302801

SUMMARY OF CLP ORGANICS ANALYSES - PHASE 1B
BERKS LANDFILL REMEDIAL INVESTIGATION
BERKS COUNTY, PENNSYLVANIA

Matrix: Groundwater Well

Semivolatile Organics

CRQL	Parameter	Sample Point GW607/18 Date Sampled: 07/01/93 Dilution Factor: 1.0 Percent Moisture: NA %			Sample Point GW607/RE Date Sampled: 07/01/93 Dilution Factor: 1.0 Percent Moisture: NA %			Sample Point GW608/18 Date Sampled: 07/01/93 Dilution Factor: 1.0 Percent Moisture: NA %			Sample Point GW610/18 Date Sampled: 07/01/93 Dilution Factor: 1.0 Percent Moisture: NA %		
		SQL	Result	Qual	SQL	Result	Qual	SQL	Result	Qual	SQL	Result	Qual
10	Acenaphthene	10	10	UL	10	10	U	10	10	U	10	10	U
10	Acenaphthylene	10	10	UL	10	10	U	10	10	U	10	10	U
10	Anthracene	10	10	UL	10	10	U	10	10	U	10	10	U
10	Benzo(a)Anthracene	10	10	UL	10	10	U	10	10	U	10	10	U
10	4-Bromophenyl-phenylether	10	10	UL	10	10	U	10	10	U	10	10	U
10	Butylbenzylphthalate	10	10	UL	10	10	U	10	10	U	10	10	U
10	Carbazole	10	10	UL	10	10	U	10	10	U	10	10	U
10	4-Chloro-3-Methylphenol	10	10	UL	10	10	U	10	10	U	10	10	U
10	4-Chloroaniline	10	10	UL	10	10	U	10	10	U	10	10	U
10	bis(2-Chloroethoxy)Methane	10	10	UL	10	10	U	10	10	U	10	10	U
10	bis(2-Chloroethyl)ether	10	10	UL	10	10	U	10	10	U	10	10	U
10	2-Chloronaphthalene	10	10	UL	10	10	U	10	10	U	10	10	U
10	2-Chlorophenol	10	10	UL	10	10	U	10	10	U	10	10	U
10	4-Chlorophenyl-phenylether	10	10	UL	10	10	U	10	10	U	10	10	U
10	Chrysene	10	10	UL	10	10	U	10	10	U	10	10	U
10	Dibenz(a,h)Anthracene	10	10	UL	10	10	U	10	10	U	10	10	U
10	Dibenzofuran	10	10	UL	10	10	U	10	10	U	10	10	U
10	1,2-Dichlorobenzene	10	10	UL	10	10	U	10	10	U	10	10	U
10	1,3-Dichlorobenzene	10	10	UL	10	10	U	10	10	U	10	10	U
10	1,4-Dichlorobenzene	10	10	UL	10	10	U	10	10	U	10	10	U
10	3,3'-Dichlorobenzidine	10	10	UL	10	10	U	10	10	U	10	10	U
10	2,4-Dichlorophenol	10	10	UL	10	10	U	10	10	U	10	10	U
10	Diethylphthalate	10	10	UL	10	10	U	10	10	U	10	10	U
10	2,4-Dimethylphenol	10	10	UL	10	10	U	10	10	U	10	10	U
10	Dimethyl Phthalate	10	10	UL	10	10	U	10	10	U	10	10	U
10	Di-n-butylphthalate	10	10	UL	10	10	U	10	10	U	10	10	U
25	4,6-Dinitro-2-methylphenol	25	25	UL	25	25	U	25	25	U	25	25	U
25	2,4-Dinitrophenol	25	25	UL	25	25	U	25	25	U	25	25	U
10	2,6-Dinitrotoluene	10	10	UL	10	10	U	10	10	U	10	10	U
10	2,6-Dinitrotoluene	10	10	UL	10	10	U	10	10	U	10	10	U
10	bis(2-Ethylhexyl)phthalate	10	10	UL	10	10	U	10	10	U	10	10	U
10	Fluorene	10	10	UL	10	10	U	10	10	U	10	10	U
10	Benzo(b)Fluorene	10	10	UL	10	10	U	10	10	U	10	10	U
10	Benzo(k)Fluorene	10	10	UL	10	10	U	10	10	U	10	10	U
10	Fluorene	10	10	UL	10	10	U	10	10	U	10	10	U
10	Hexachlorobenzene	10	10	UL	10	10	U	10	10	U	10	10	U
10	Hexachlorobutadiene	10	10	UL	10	10	U	10	10	U	10	10	U
10	Hexachlorocyclopentadiene	10	10	UL	10	10	U	10	10	U	10	10	U
10	Hexachloroethene	10	10	UL	10	10	U	10	10	U	10	10	U

Notes:

All units are ug/l
Sample Quantitation Limit (SQL) is CRQL multiplied by the dilution factor and adjusted for percent moisture as appropriate.
The Qual column indicates the qualifier applied to the result following data validation (see below).

Qualifiers:
A - Acceptable (Quantitative) Data K - Biased High (Semi-quantitative) Data UJ - Not Detected/Estimated Data N - Tentative Identification
J - Estimated (Semi-quantitative) Data L - Biased Low (Semi-quantitative) Data U - Not Detected/Quantitative Data R - Unusable Data
B - Present in associated blanks (Semi-quantitative Data) UL - Not Detected/Biased Low UL - Not Detected/Biased Low Q - Not Applicable

AR302802

Matrix: Groundwater Well

Semivolatile Organics

CRQL	Parameter	Sample Point GUG07/1B			Sample Point GUG07/BE			Sample Point GUG08/1B			Sample Point GUG10/1B		
		SOL	Result	Qual	SOL	Result	Qual	SOL	Result	Qual	SOL	Result	Qual
10	Indeno(1,2,3-cd)Pyrene	10	10	UL	10	10	U	10	10	U	10	10	U
10	Isophorone	10	10	UL	10	10	U	10	10	U	10	10	U
10	2-Methylnaphthalene	10	10	UL	10	10	U	10	10	U	10	10	U
10	2-Methylphenol	10	10	UL	10	10	U	10	10	U	10	10	U
10	4-Methylphenol	10	10	UL	10	10	U	10	10	U	10	10	U
10	Naphthalene	10	10	UL	10	10	U	10	10	U	10	10	U
25	2-Nitroaniline	25	25	UL	25	25	U	25	25	U	25	25	U
25	3-Nitroaniline	25	25	UL	25	25	U	25	25	U	25	25	U
25	4-Nitroaniline	25	25	UL	25	25	U	25	25	U	25	25	U
10	Nitrobenzene	10	10	UL	10	10	U	10	10	U	10	10	U
10	2-Nitrophenol	10	10	UL	10	10	U	10	10	U	10	10	U
25	4-Nitrophenol	25	25	UL	25	25	U	25	25	U	25	25	U
10	N-Nitroso-di-n-propylamine	10	10	UL	10	10	U	10	10	U	10	10	U
10	N-Nitrosodiphenylamine	10	10	UL	10	10	U	10	10	U	10	10	U
10	2,2'-Oxybis(1-Chloropropane)	10	10	UL	10	10	U	10	10	U	10	10	U
25	Pentachlorophenol	25	25	UL	25	25	U	25	25	U	25	25	U
10	Benzo(g,h,i)perylene	10	10	UL	10	10	U	10	10	U	10	10	U
10	Phenanthrene	10	10	UL	10	10	U	10	10	U	10	10	U
10	Pyrene	10	10	UL	10	10	U	10	10	U	10	10	U
10	Benzo(a)Pyrene	10	10	UL	10	10	U	10	10	U	10	10	U
10	1,2,4-Trichlorobenzene	10	10	UL	10	10	U	10	10	U	10	10	U
25	2,4,5-Trichlorophenol	25	25	UL	25	25	U	25	25	U	25	25	U
10	2,4,6-Trichlorophenol	10	10	UL	10	10	U	10	10	U	10	10	U

Notes:
All units are ug/l
Sample Quantitation Limit (SOL) is CRQL multiplied by the dilution factor and adjusted for percent moisture as appropriate.
The qual column indicates the qualifier applied to the result following data validation (see below).

Qualifiers:.....
A - Acceptable (Quantitative) Data K - Biased High (Semi-quantitative) Data
J - Estimated (Semi-quantitative) Data L - Biased Low (Semi-quantitative) Data
B - Present in associated blanks (Semi-quantitative Data)
UL - Not Detected/Biased Low
UJ - Not Detected/Estimated Data
U - Not Detected/Quantitative Data
H - Tentative Identification
R - Unusable Data
Q - Not Applicable

AR302803

SUMMARY OF CLP ORGANIC ANALYSES - PHASE 1B
BERKS LANDFILL REMEDIAL INVESTIGATION
BERKS COUNTY, PENNSYLVANIA

Matrix: Groundwater Well

Semivolatile Organics

CROL	Parameter	Sample Point GMG11/1B			Sample Point GMG11/RE			Sample Point GMG11/FD			Sample Point GMG11/FR		
		SQL	Result	Qual	SQL	Result	Qual	SQL	Result	Qual	SQL	Result	Qual
10	Acenaphthene	10	10	U	10	10	U	10	10	U	10	10	U
10	Acenaphthylene	10	10	U	10	10	U	10	10	U	10	10	U
10	Anthracene	10	10	U	10	10	U	10	10	U	10	10	U
10	Benzo(a)Anthracene	10	10	U	10	10	U	10	10	U	10	10	U
10	4-Bromophenyl-phenylether	10	10	U	10	10	U	10	10	U	10	10	U
10	Butylbenzylphthalate	10	10	U	10	10	U	10	10	U	10	10	U
10	Carbazole	10	10	U	10	10	U	10	10	U	10	10	U
10	4-Chloro-3-Methylphenol	10	10	U	10	10	U	10	10	U	10	10	U
10	4-Chloroaniline	10	10	U	10	10	U	10	10	U	10	10	U
10	bis(2-Chloroethoxy)Methane	10	10	U	10	10	U	10	10	U	10	10	U
10	bis(2-Chloroethyl)ether	10	10	U	10	10	U	10	10	U	10	10	U
10	2-Chloronaphthalene	10	10	U	10	10	U	10	10	U	10	10	U
10	2-Chlorophenol	10	10	U	10	10	U	10	10	U	10	10	U
10	4-Chlorophenyl-phenylether	10	10	U	10	10	U	10	10	U	10	10	U
10	Chrysene	10	10	U	10	10	U	10	10	U	10	10	U
10	Dibenz(a,h)Anthracene	10	10	U	10	10	U	10	10	U	10	10	U
10	Dibenzofuran	10	10	U	10	10	U	10	10	U	10	10	U
10	1,2-Dichlorobenzene	10	2	J	10	2	J	10	2	J	10	2	J
10	1,3-Dichlorobenzene	10	9	J	10	10	J	10	10	J	10	10	J
10	1,4-Dichlorobenzene	10	10	A	10	10	A	10	10	A	10	10	A
10	3,3'-Dichlorobenzidine	10	10	U	10	10	U	10	10	U	10	10	U
10	2,4-Dichlorophenol	10	10	U	10	10	U	10	10	U	10	10	U
10	Diethylphthalate	10	7	B	10	7	B	10	6	B	10	7	B
10	2,4-Dimethylphenol	10	10	U	10	10	U	10	10	U	10	10	U
10	Dimethyl Phthalate	10	10	U	10	10	U	10	10	U	10	10	U
10	Di-n-butylphthalate	10	1	B	10	1	B	10	1	B	10	1	B
25	4,6-Dinitro-2-methylphenol	25	25	U	25	25	U	25	25	U	25	25	U
25	2,4-Dinitrophenol	25	25	U	25	25	U	25	25	U	25	25	U
10	2,4-Dinitrotoluene	10	10	U	10	10	U	10	10	U	10	10	U
10	2,6-Dinitrotoluene	10	10	U	10	10	U	10	10	U	10	10	U
10	Di-n-Octyl Phthalate	10	10	U	10	10	U	10	10	U	10	10	U
10	bis(2-Ethylhexyl)Phthalate	10	2	B	10	2	B	10	3	B	10	3	B
10	Fluoranthene	10	10	U	10	10	U	10	10	U	10	10	U
10	Benzo(b)Fluoranthene	10	10	U	10	10	U	10	10	U	10	10	U
10	Benzo(k)Fluoranthene	10	10	U	10	10	U	10	10	U	10	10	U
10	Fluorene	10	10	U	10	10	U	10	10	U	10	10	U
10	Hexachlorobenzene	10	10	U	10	10	U	10	10	U	10	10	U
10	Hexachlorobutadiene	10	10	U	10	10	U	10	10	U	10	10	U
10	Hexachlorocyclopentadiene	10	10	U	10	10	U	10	10	U	10	10	U
10	Hexachloroethane	10	6	J	10	6	J	10	6	J	10	6	J

Notes:
All units are ug/l
Sample Quantitation Limit (SQL) is CROL multiplied by the dilution factor and adjusted for percent moisture as appropriate.
The Qual column indicates the qualifier applied to the result following data validation (see below).

Qualifiers:
A - Acceptable (Quantitative) Data
K - Biased High (Semi-quantitative) Data
L - Biased Low (Semi-quantitative) Data
J - Estimated (Semi-quantitative) Data
B - Present in associated blanks (Semi-quantitative) Data
UJ - Not Detected/Estimated Data
U - Not Detected/Quantitative Data
UL - Not Detected/Biased Low
N - Tentative Identification
R - Unusable Data
Q - Not Applicable

AR302804

SUMMARY OF CLP ORGANIC ANALYSES - PHASE 1B
BERKS LANDFILL REMEDIAL INVESTIGATION
BERKS COUNTY, PENNSYLVANIA

Matrix: Groundwater Well

Semivolatile Organics

CRQL	Parameter	Sample Point GUG11/1b			Sample Point GUG11/RE			Sample Point GUG11/FO			Sample Point GUG11/FR		
		SQL	Result	Qual	SQL	Result	Qual	SQL	Result	Qual	SQL	Result	Qual
10	Indeno(1,2,3-cd)Pyrene	10	10	UJ	10	10	UJ	10	10	UJ	10	10	UJ
10	Isophorone	10	10	U	10	10	U	10	10	U	10	10	U
10	2-Methylnaphthalene	10	10	U	10	10	U	10	10	U	10	10	U
10	2-Methylphenol	10	10	U	10	10	U	10	10	U	10	10	U
10	4-Methylphenol	10	10	U	10	10	U	10	10	U	10	10	U
10	Naphthalene	10	10	U	10	10	U	10	10	U	10	10	U
25	2-Nitroaniline	25	25	U	25	25	U	25	25	U	25	25	U
25	3-Nitroaniline	25	1	J	25	25	U	25	25	U	25	25	U
25	4-Nitroaniline	25	7	J	25	25	U	25	25	U	25	25	U
10	Nitrobenzene	10	10	U	10	10	U	10	10	U	10	10	U
10	2-Nitrophenol	10	10	U	10	10	U	10	10	U	10	10	U
25	4-Nitrophenol	25	25	U	25	25	U	25	25	U	25	25	U
10	N-Nitroso-di-n-propylamine	10	1	J	10	10	U	10	10	U	10	10	U
10	N-Nitrosodiphenylamine	10	10	U	10	10	U	10	10	U	10	10	U
10	2,2'-Oxybis(1-Chloropropane)	10	10	U	10	10	U	10	10	U	10	10	U
25	Pentachlorophenol	25	25	U	25	25	U	25	25	U	25	25	U
10	Benzo(g,h,i)Perylene	10	10	UJ	10	10	UJ	10	10	UJ	10	10	UJ
10	Phenanthrene	10	10	U	10	10	U	10	10	U	10	10	U
10	Phenol	10	10	U	10	10	U	10	10	U	10	10	U
10	Pyrene	10	10	U	10	10	U	10	10	U	10	10	U
10	Benzo(a)Pyrene	10	10	UJ	10	10	UJ	10	10	UJ	10	10	UJ
10	1,2,4-Trichlorobenzene	10	10	U	10	10	U	10	10	U	10	10	U
25	2,4,5-Trichlorophenol	25	25	U	25	25	U	25	25	U	25	25	U
10	2,4,6-Trichlorophenol	10	10	U	10	10	U	10	10	U	10	10	U

Notes:
All units are ug/l
Sample Quantitation Limit (SQL) is CRQL multiplied by the dilution factor and adjusted for percent moisture as appropriate.
The qual column indicates the qualifier applied to the result following data validation (see below).

Qualifiers:
A - Acceptable (Quantitative) Data
J - Estimated (Semi-quantitative) Data
B - Present in associated blanks (Semi-quantitative Data)
K - Biased High (Semi-quantitative) Data
L - Biased Low (Semi-quantitative) Data
UJ - Not Detected/Estimated Data
UL - Not Detected/Quantitative Data
M - Tentative Identification
N - Unusable Data
Q - Not Applicable
R - Not Detected/Quantitative Data
UL - Not Detected/Biased Low

AR302805

Matrix: Groundwater Well

Semivolatile Organics

CRQL	Parameter	Sample Point GMG12/18			Sample Point GMG12/18			Sample Point GMG12/18			Sample Point GMG13/18			Sample Point GMG03/18		
		SQL	Result	Qual	SQL	Result	Qual	SQL	Result	Qual	SQL	Result	Qual	SQL	Result	Qual
10	Acenaphthene	10	10	U	10	10	U	10	10	U	10	10	U	10	10	U
10	Acenaphthylene	10	10	U	10	10	U	10	10	U	10	10	U	10	10	U
10	Anthracene	10	10	U	10	10	U	10	10	U	10	10	U	10	10	U
10	Benzo(a)Anthracene	10	10	U	10	10	U	10	10	U	10	10	U	10	10	U
10	4-Bromophenyl-phenylether	10	10	U	10	10	U	10	10	U	10	10	U	10	10	U
10	Butylbenzylphthalate	10	10	U	10	10	U	10	10	U	10	10	U	10	10	U
10	Carbazole	10	10	U	10	10	U	10	10	U	10	10	U	10	10	U
10	4-Chloro-3-Methylphenol	10	10	U	10	10	U	10	10	U	10	10	U	10	10	U
10	4-Chloroaniline	10	10	U	10	10	U	10	10	U	10	10	U	10	10	U
10	bis(2-Chloroethoxy)Methane	10	10	U	10	10	U	10	10	U	10	10	U	10	10	U
10	bis(2-Chloroethyl)ether	10	10	U	10	10	U	10	10	U	10	10	U	10	10	U
10	2-Chloronaphthalene	10	10	U	10	10	U	10	10	U	10	10	U	10	10	U
10	2-Chlorophenol	10	10	U	10	10	U	10	10	U	10	10	U	10	10	U
10	4-Chlorophenyl-phenylether	10	10	U	10	10	U	10	10	U	10	10	U	10	10	U
10	Chrysene	10	10	U	10	10	U	10	10	U	10	10	U	10	10	U
10	Dibenz(a,h)Anthracene	10	10	U	10	10	U	10	10	U	10	10	U	10	10	U
10	Dibenzofuran	10	10	U	10	10	U	10	10	U	10	10	U	10	10	U
10	1,2-Dichlorobenzene	10	10	U	10	10	U	10	10	U	10	10	U	10	10	U
10	1,3-Dichlorobenzene	10	10	U	10	10	U	10	10	U	10	10	U	10	10	U
10	1,4-Dichlorobenzene	10	10	U	10	10	U	10	10	U	10	10	U	10	10	U
10	3,3'-Dichlorobenzidine	10	2	J	10	2	J	10	2	J	10	2	J	10	2	J
10	2,4-Dichlorophenol	10	10	U	10	10	U	10	10	U	10	10	U	10	10	U
10	Diethylphthalate	10	10	U	10	10	U	10	10	U	10	10	U	10	10	U
10	2,4-Dimethylphenol	10	10	U	10	10	U	10	10	U	10	10	U	10	10	U
10	Dimethyl Phthalate	10	10	U	10	10	U	10	10	U	10	10	U	10	10	U
10	Di-n-butylphthalate	10	10	U	10	10	U	10	10	U	10	10	U	10	10	U
25	4,6-Dinitro-2-methylphenol	25	25	U	25	25	U	25	25	U	25	25	U	25	25	U
25	2,4-Dinitrophenol	25	25	U	25	25	U	25	25	U	25	25	U	25	25	U
10	2,4-Dinitrotoluene	10	10	U	10	10	U	10	10	U	10	10	U	10	10	U
10	2,6-Dinitrotoluene	10	10	U	10	10	U	10	10	U	10	10	U	10	10	U
10	Di-n-Octyl Phthalate	10	10	U	10	10	U	10	10	U	10	10	U	10	10	U
10	bis(2-Ethylhexyl)phthalate	10	10	U	10	10	U	10	10	U	10	10	U	10	10	U
10	Fluoranthene	10	10	U	10	10	U	10	10	U	10	10	U	10	10	U
10	Benzo(b)Fluoranthene	10	10	U	10	10	U	10	10	U	10	10	U	10	10	U
10	Benzo(k)Fluoranthene	10	10	U	10	10	U	10	10	U	10	10	U	10	10	U
10	Fluorene	10	10	U	10	10	U	10	10	U	10	10	U	10	10	U
10	Hexachlorobenzene	10	10	U	10	10	U	10	10	U	10	10	U	10	10	U
10	Hexachlorobutadiene	10	10	U	10	10	U	10	10	U	10	10	U	10	10	U
10	Hexachlorocyclopentadiene	10	10	U	10	10	U	10	10	U	10	10	U	10	10	U
10	Hexachloroethene	10	10	U	10	10	U	10	10	U	10	10	U	10	10	U

Notes:

All units are ug/l
Sample Quantitation Limit (SQL) is CRQL multiplied by the dilution factor and adjusted for percent moisture as appropriate.
The Qual column indicates the qualifier applied to the result following data validation (see below).

Qualifiers:.....
A - Acceptable (Quantitative) Data
J - Estimated (Semi-quantitative) Data
B - Present in associated blanks (Semi-quantitative Data)
K - Baised High (Semi-quantitative) Data
L - Baised Low (Semi-quantitative) Data
UJ - Not Detected/Estimated Data
U - Not Detected/Quantitative Data
UL - Not Detected/Baised Low
N - Tentative Identification
R - Unusable Data
Q - Not Applicable

AR302806

Matrix: Groundwater Well

Semivolatile Organics

CRQL	Parameter	Sample Point GUG12/1B			Sample Point GUG12/1B			Sample Point GUG12/1B			Sample Point GUG13/1B			Sample Point GUG03/1B		
		SQL	Result	Qual	SQL	Result	Qual	SQL	Result	Qual	SQL	Result	Qual	SQL	Result	Qual
10	Indeno(1,2,3-cd)Pyrene	10	10	U	10	10	U	10	10	U	10	10	10	10	U	U
10	Isophorone	10	10	U	10	10	U	10	10	U	10	10	10	10	U	U
10	2-Methylnaphthalene	10	10	U	10	10	U	10	10	U	10	10	10	10	U	U
10	2-Methylphenol	10	10	U	10	10	U	10	10	U	10	10	10	10	U	U
10	4-Methylphenol	10	10	U	10	10	U	10	10	U	10	10	10	10	U	U
10	Naphthalene	10	10	U	10	10	U	10	10	U	10	10	10	10	U	U
25	2-Nitroaniline	25	25	U	25	25	U	25	25	U	25	25	25	25	U	U
25	3-Nitroaniline	25	25	U	25	25	U	25	25	U	25	25	25	25	U	U
25	4-Nitroaniline	25	25	U	25	25	U	25	25	U	25	25	25	25	U	U
10	Nitrobenzene	10	10	U	10	10	U	10	10	U	10	10	10	10	U	U
10	2-Nitrophenol	10	10	U	10	10	U	10	10	U	10	10	10	10	U	U
25	4-Nitrophenol	25	25	U	25	25	U	25	25	U	25	25	25	25	U	U
10	N-Nitroso-di-n-propylamine	10	10	U	10	10	U	10	10	U	10	10	10	10	U	U
10	N-Nitrosodiphenylamine	10	10	U	10	10	U	10	10	U	10	10	10	10	U	U
10	2,2'-Oxybis(1-Chloropropane)	10	10	U	10	10	U	10	10	U	10	10	10	10	U	U
25	Pentachlorophenol	25	25	U	25	25	U	25	25	U	25	25	25	25	U	U
10	Benzo(g,h,i)Perylene	10	10	U	10	10	U	10	10	U	10	10	10	10	U	U
10	Phenanthrene	10	10	U	10	10	U	10	10	U	10	10	10	10	U	U
10	Phenol	10	10	U	10	10	U	10	10	U	10	10	10	10	U	U
10	Pyrene	10	10	U	10	10	U	10	10	U	10	10	10	10	U	U
10	Benzo(a)Pyrene	10	10	U	10	10	U	10	10	U	10	10	10	10	U	U
10	1,2,4-Trichlorobenzene	10	10	U	10	10	U	10	10	U	10	10	10	10	U	U
25	2,4,5-Trichlorophenol	25	25	U	25	25	U	25	25	U	25	25	25	25	U	U
10	2,6-Dichlorophenol	10	10	U	10	10	U	10	10	U	10	10	10	10	U	U

Notes:
All units are ug/l
Sample Quantitation Limit (SQL) is CRQL multiplied by the dilution factor and adjusted for percent moisture as appropriate.
The qual column indicates the qualifier applied to the result following data validation (see below).

Qualifiers:
A - Acceptable (Quantitative) Data
J - Estimated (Semi-quantitative) Data
K - Biased High (Semi-quantitative) Data
L - Biased Low (Semi-quantitative) Data
M - Present in associated blanks (Semi-quantitative Data)
N - Not Detected/Estimated Data
OJ - Not Detected/Quantitative Data
P - Unusable Data
Q - Not Applicable
R - Unusable Data
UL - Not Detected/Biased Low

AR302807

SUMMARY OF CLP ORGANICS ANALYSES - PHASE 1B
BERKS LANDFILL REMEDIAL INVESTIGATION
BERKS COUNTY, PENNSYLVANIA

Matrix: Groundwater Well

Semivolatile Organics

CROL	Parameter	Sample Point GM06/1B			Sample Point GM11/1B			Sample Point GM140/1B			Sample Point GM145/1B		
		SQL	Result	Qual	SQL	Result	Qual	SQL	Result	Qual	SQL	Result	Qual
10	Acenaphthene	10	10	R	10	10	U	10	10	U	10	10	U
10	Acenaphthylene	10	10	R	10	10	U	10	10	U	10	10	U
10	Anthracene	10	10	R	10	10	U	10	10	U	10	10	U
10	Benzo(a)Anthracene	10	10	R	10	10	U	10	10	U	10	10	U
10	4-Bromophenyl-phenylether	10	10	R	10	10	U	10	10	U	10	10	U
10	Butylbenzylphthalate	10	10	R	10	10	U	10	10	U	10	10	U
10	Carbazole	10	10	R	10	10	U	10	10	U	10	10	U
10	4-Chloro-3-Methylphenol	10	10	R	10	10	U	10	10	U	10	10	U
10	4-Chloroaniline	10	10	R	10	10	U	10	10	U	10	10	U
10	bis(2-Chloroethoxy)Methane	10	10	R	10	10	U	10	10	U	10	10	U
10	bis(2-Chloroethyl)ether	10	10	R	10	10	U	10	10	U	10	10	U
10	2-Chloronaphthalene	10	10	R	10	10	U	10	10	U	10	10	U
10	2-Chlorophenol	10	10	R	10	10	U	10	10	U	10	10	U
10	4-Chlorophenyl-phenylether	10	10	R	10	10	U	10	10	U	10	10	U
10	Chrysene	10	10	R	10	10	U	10	10	U	10	10	U
10	Dibenz(a,h)Anthracene	10	10	R	10	10	U	10	10	U	10	10	U
10	Dibenzofuran	10	10	R	10	10	U	10	10	U	10	10	U
10	1,2-Dichlorobenzene	10	10	R	10	10	U	10	10	U	10	10	U
10	1,3-Dichlorobenzene	10	10	R	10	10	U	10	10	U	10	10	U
10	1,4-Dichlorobenzene	10	10	R	10	10	U	10	10	U	10	10	U
10	3,3'-Dichlorobenzidine	10	10	R	10	10	U	10	10	U	10	10	U
10	2,4-Dichlorophenol	10	10	R	10	10	U	10	10	U	10	10	U
10	Diethylphthalate	10	10	R	10	10	U	10	10	U	10	10	U
10	2,4-Dimethylphenol	10	10	R	10	10	U	10	10	U	10	10	U
10	Dimethyl Phthalate	10	10	R	10	10	U	10	10	U	10	10	U
10	Di-n-butylphthalate	10	10	R	10	10	U	10	10	U	10	10	U
25	4,6-Dinitro-2-methylphenol	25	25	R	25	25	U	25	25	U	25	25	U
25	2,4-Dinitrophenol	25	25	R	25	25	U	25	25	U	25	25	U
10	2,4-Dinitrotoluene	10	10	R	10	10	U	10	10	U	10	10	U
10	2,6-Dinitrotoluene	10	10	R	10	10	U	10	10	U	10	10	U
10	Di-n-Octyl Phthalate	10	10	R	10	10	U	10	10	U	10	10	U
10	bis(2-Ethylhexyl)Phthalate	10	10	R	10	10	U	10	10	U	10	10	U
10	Fluoranthene	10	10	R	10	10	U	10	10	U	10	10	U
10	Benzo(b)Fluoranthene	10	10	R	10	10	U	10	10	U	10	10	U
10	Benzo(k)Fluoranthene	10	10	R	10	10	U	10	10	U	10	10	U
10	Fluorene	10	10	R	10	10	U	10	10	U	10	10	U
10	Hexachlorobenzene	10	10	R	10	10	U	10	10	U	10	10	U
10	Hexachlorobutadiene	10	10	R	10	10	U	10	10	U	10	10	U
10	Hexachlorocyclopentadiene	10	10	R	10	10	U	10	10	U	10	10	U
10	Hexachloroethane	10	10	R	10	10	U	10	10	U	10	10	U

Notes:

All units are ug/l
Sample Quantitation Limit (SQL) is CROL multiplied by the dilution factor and adjusted for percent moisture as appropriate.
The qual column indicates the qualifier applied to the result following data validation (see below).

Qualifiers.....
A - Acceptable (Quantitative) Data
K - Biased High (Semi-quantitative) Data
L - Biased Low (Semi-quantitative) Data
B - Present in associated blanks (Semi-quantitative Data)
UJ - Not Detected/Estimated Data
U - Not Detected/Quantitative Data
UL - Not Detected/Biased Low
N - Tentative Identification
R - Unusable Data
Q - Not Applicable

Matrix: Groundwater Well

Semivolatile Organics

CRQL	Parameter	Sample Point GIM06/1B			Sample Point GIM11/1B			Sample Point GIM140/1B			Sample Point GIM145/1B		
		SQL	Result	Qual	SQL	Result	Qual	SQL	Result	Qual	SQL	Result	Qual
10	Indeno(1,2,3-cd)Pyrene	10	10	R	10	10	U	10	10	U	10	10	U
10	Isophorone	10	10	R	10	10	U	10	10	U	10	10	U
10	2-Methylnaphthalene	10	10	R	10	10	U	10	10	U	10	10	U
10	2-Methylphenol	10	10	U	10	10	U	10	10	U	10	10	U
10	4-Methylphenol	10	10	U	10	10	U	10	10	U	10	10	U
10	Naphthalene	10	2	L	10	10	U	10	10	U	10	10	U
25	2-Nitroaniline	25	25	R	25	25	U	25	25	U	25	25	U
25	3-Nitroaniline	25	25	R	25	25	U	25	25	U	25	25	U
25	4-Nitroaniline	25	25	R	25	25	U	25	25	U	25	25	U
10	Nitrobenzene	10	10	U	10	10	U	10	10	U	10	10	U
10	2-Nitrophenol	10	10	U	10	10	U	10	10	U	10	10	U
25	4-Nitrophenol	25	25	U	25	25	U	25	25	U	25	25	U
10	N-Nitroso-di-n-propylamine	10	10	R	10	10	U	10	10	U	10	10	U
10	N-Nitrosodiphenylamine	10	10	R	10	10	U	10	10	U	10	10	U
10	2,2'-Oxybis(1-Chloropropane)	10	10	R	10	10	U	10	10	U	10	10	U
25	Pentachlorophenol	25	25	U	25	25	U	25	25	U	25	25	U
10	Benzo(g,h,i)Perylene	10	10	R	10	10	U	10	10	U	10	10	U
10	Phenanthrene	10	10	U	10	10	U	10	10	U	10	10	U
10	Phenol	10	10	U	10	10	U	10	10	U	10	10	U
10	Pyrene	10	10	R	10	10	U	10	10	U	10	10	U
10	Benzo(a)Pyrene	10	10	R	10	10	U	10	10	U	10	10	U
10	1,2,4-Trichlorobenzene	10	10	R	10	10	U	10	10	U	10	10	U
25	2,4,5-Trichlorophenol	25	25	U	25	25	U	25	25	U	25	25	U
10	2,4,6-Trichlorophenol	10	10	U	10	10	U	10	10	U	10	10	U

Notes:
All units are ug/l
Sample Quantitation Limit (SQL) is CRQL multiplied by the dilution factor and adjusted for percent moisture as appropriate.
The Qual column indicates the qualifier applied to the result following data validation (see below).

Qualifiers:.....
A - Acceptable (Quantitative) Data K - Biased High (Semi-quantitative) Data UJ - Not Detected/Estimated Data M - Tentative Identification
J - Estimated (Semi-quantitative) Data L - Biased Low (Semi-quantitative) Data U - Not Detected/Quantitative Data R - Unusable Data
B - Present in associated blanks (Semi-quantitative Data) UL - Not Detected/Biased Low UL - Not Detectable

AR302809

SUMMARY OF CLP ORGANICS ANALYSES - PHASE 1B
BERKS LANDFILL REMEDIAL INVESTIGATION
BERKS COUNTY, PENNSYLVANIA

Matrix: Groundwater Well

Semivolatile Organics

CROL	Parameter	Sample Point GM14S/RE			Sample Point GM15D/1B			Sample Point GM15S/1B			Sample Point GM16/1B		
		SQL	Result	Qual	SQL	Result	Qual	SQL	Result	Qual	SQL	Result	Qual
10	Acenaphthene	10	10	U	10	10	U	10	10	U	10	10	U
10	Acenaphthylene	10	10	U	10	10	U	10	10	U	10	10	U
10	Anthracene	10	10	U	10	10	U	10	10	U	10	10	U
10	Benzo(a)Anthracene	10	10	U	10	10	U	10	10	U	10	10	U
10	4-Bromophenyl-phenylether	10	10	U	10	10	U	10	10	U	10	10	U
10	Butylbenzylphthalate	10	10	U	10	10	U	10	10	U	10	10	U
10	Carbazole	10	10	U	10	10	U	10	10	U	10	10	U
10	4-Chloro-3-Methylphenol	10	10	U	10	10	U	10	10	U	10	10	U
10	4-Chloroaniline	10	10	U	10	10	U	10	10	U	10	10	U
10	bis(2-Chloroethoxy)Methane	10	10	U	10	10	U	10	10	U	10	10	U
10	bis(2-Chloroethyl)ether	10	10	U	10	10	U	10	10	U	10	10	U
10	2-Chloronaphthalene	10	10	U	10	10	U	10	10	U	10	10	U
10	2-Chlorophenol	10	10	U	10	10	U	10	10	U	10	10	U
10	4-Chlorophenyl-phenylether	10	10	U	10	10	U	10	10	U	10	10	U
10	Chrysene	10	10	U	10	10	U	10	10	U	10	10	U
10	Dibenz(a,h)Anthracene	10	10	U	10	10	U	10	10	U	10	10	U
10	Dibenzofuran	10	10	U	10	10	U	10	10	U	10	10	U
10	1,2-Dichlorobenzene	10	10	U	10	10	U	10	10	U	10	10	U
10	1,3-Dichlorobenzene	10	10	U	10	10	U	10	10	U	10	10	U
10	1,4-Dichlorobenzene	10	10	U	10	10	U	10	10	U	10	10	U
10	3,3'-Dichlorobenzidine	10	7	J	10	10	U	10	10	U	10	3	J
10	2,4-Dichlorophenol	10	10	U	10	10	U	10	10	U	10	10	U
10	Diethylphthalate	10	1	B	10	10	U	10	10	U	10	13	B
10	2,4-Dimethylphenol	10	10	U	10	10	U	10	10	U	10	4	J
10	Dimethyl Phthalate	10	10	U	10	10	U	10	10	U	10	10	U
10	Di-n-butylphthalate	10	10	U	10	10	U	10	10	U	10	10	U
25	4,6-Dinitro-2-methylphenol	25	25	U	25	25	U	25	25	U	25	25	U
25	2,4-Dinitrophenol	25	25	U	25	25	U	25	25	U	25	25	U
10	2,4-Dinitrotoluene	10	10	U	10	10	U	10	10	U	10	10	U
10	2,6-Dinitrotoluene	10	10	U	10	10	U	10	10	U	10	10	U
10	Di-n-Octyl Phthalate	10	10	U	10	10	U	10	10	U	10	10	U
10	bis(2-Ethylhexyl)Phthalate	10	10	U	10	10	U	10	10	U	10	2	B
10	Fluoranthene	10	10	U	10	10	U	10	10	U	10	10	U
10	Benzo(b)Fluoranthene	10	10	U	10	10	U	10	10	U	10	10	U
10	Benzo(k)Fluoranthene	10	10	U	10	10	U	10	10	U	10	10	U
10	Fluorene	10	10	U	10	10	U	10	10	U	10	10	U
10	Hexachlorobenzene	10	10	U	10	10	U	10	10	U	10	10	U
10	Hexachlorobutadiene	10	10	U	10	10	U	10	10	U	10	10	U
10	Hexachlorocyclopentadiene	10	10	U	10	10	U	10	10	U	10	10	U
10	Hexachloroethane	10	10	U	10	10	U	10	10	U	10	10	U

Notes:
All units are ug/l
Sample Quantitation Limit (SQL) is CRQL multiplied by the dilution factor and adjusted for percent moisture as appropriate.
The Qual column indicates the qualifier applied to the result following data validation (see below).

Qualifiers:.....
A - Acceptable (Quantitative) Data
J - Estimated (Semi-quantitative) Data
B - Present in associated blanks (Semi-quantitative Data)
K - Biased High (Semi-quantitative) Data
L - Biased Low (Semi-quantitative) Data
UJ - Not Detected/Estimated Data
U - Not Detected/Quantitative Data
UL - Not Detected/Biased Low
M - Tentative Identification
R - Unusable Data
Q - Not Applicable

AR302810

Matrix: Groundwater Well

Sealvolatile Organics

CRQL	Parameter	Sample Point GUM14S/RE			Sample Point GUM15D/1B			Sample Point GUM15S/1B			Sample Point GUM16/1B		
		SQL	Result	Qual	SQL	Result	Qual	SQL	Result	Qual	SQL	Result	Qual
10	Indeno(1,2,3-cd)Pyrene	10	10	UJ	10	10	U	10	10	U	10	10	U
10	Isophorone	10	10	U	10	10	U	10	10	U	10	10	U
10	2-Methylnaphthalene	10	10	U	10	10	U	10	10	U	10	10	U
10	2-Methylphenol	10	10	U	10	10	U	10	10	U	10	10	U
10	4-Methylphenol	10	10	U	10	10	U	10	10	U	10	10	U
10	Naphthalene	10	5	J	10	10	U	10	10	U	10	10	U
25	2-Nitroaniline	25	25	U	25	25	U	25	25	U	25	25	U
25	3-Nitroaniline	25	25	U	25	25	U	25	25	U	25	25	U
25	4-Nitroaniline	25	25	U	25	25	U	25	25	U	25	25	U
10	Nitrobenzene	10	10	U	10	10	U	10	10	U	10	10	U
10	2-Nitrophenol	10	10	U	10	10	U	10	10	U	10	10	U
25	4-Nitrophenol	25	25	U	25	25	U	25	25	U	25	25	U
10	N-Nitroso-di-n-propylamine	10	10	U	10	10	U	10	10	U	10	10	U
10	N-Nitrosodiphenylamine	10	10	U	10	10	U	10	10	U	10	10	U
10	2,2'-Oxybis(1-Chloropropane)	10	10	U	10	10	U	10	10	U	10	10	U
25	Pentachlorophenol	25	25	UJ	25	25	U	25	25	U	25	25	U
10	Benzot(g,h,i)perylene	10	10	U	10	10	U	10	10	U	10	10	U
10	Phenanthrene	10	10	U	10	10	U	10	10	U	10	10	U
10	Phenol	10	10	U	10	2	B	10	10	U	10	10	U
10	Pyrene	10	10	U	10	10	U	10	10	U	10	10	U
10	Benzot(a)Pyrene	10	10	UJ	10	10	U	10	10	U	10	10	U
10	1,2,4-Trichlorobenzene	10	10	U	10	10	U	10	10	U	10	10	U
25	2,4,5-Trichlorophenol	25	25	U	25	25	U	25	25	U	25	25	U
10	2,4,6-Trichlorophenol	10	10	U	10	10	U	10	10	U	10	10	U

Notes:
All units are ug/l
Sample Quantitation Limit (SQL) is CRQL multiplied by the dilution factor and adjusted for percent moisture as appropriate.
The Qual column indicates the qualifier applied to the result following data validation (see below).

Qualifiers:.....
A - Acceptable (Quantitative) Data K - Biased High (Semi-quantitative) Data UJ - Not Detected/Estimated Data M - Tentative Identification
J - Estimated (Semi-quantitative) Data L - Biased Low (Semi-quantitative) Data UL - Not Detected/Quantitative Data R - Unusable Data
B - Present in associated blanks (Semi-quantitative Data) UL - Not Detected/Biased Low Q - Not Applicable

AR302811

Matrix: Groundwater Well

Semivolatile Organics

CROL	Parameter	Sample Point GM17/1B		Sample Point GM19S/1B		Sample Point GMR190/1B		Sample Point Date Sampled: Dilution Factor: Percent Moisture: %
		SQL	Result	SQL	Result	SQL	Result	
10	Acenaphthene	10	10	10	10	10	10	
10	Acenaphthylene	10	10	10	10	10	10	
10	Anthracene	10	10	10	10	10	10	
10	Benzo(a)Anthracene	10	10	10	10	10	10	
10	4-Bromophenyl-phthalate	10	10	10	10	10	10	
10	Butylbenzylphthalate	10	10	10	10	10	10	
10	Carbazole	10	10	10	10	10	10	
10	4-Chloro-3-Methylphenol	10	10	10	10	10	10	
10	4-Chloroaniline	10	10	10	10	10	10	
10	bis(2-Chloroethoxy)Methane	10	10	10	10	10	10	
10	bis(2-Chloroethyl)ether	10	10	10	10	10	10	
10	2-Chloronaphthalene	10	10	10	10	10	10	
10	2-Chlorophenol	10	10	10	10	10	10	
10	4-Chlorophenyl-phnylether	10	10	10	10	10	10	
10	Chrysene	10	10	10	10	10	10	
10	Dibenz(a,h)Anthracene	10	10	10	10	10	10	
10	Dibenzofuran	10	10	10	10	10	10	
10	1,2-Dichlorobenzene	10	10	10	10	10	10	
10	1,3-Dichlorobenzene	10	10	10	10	10	10	
10	1,4-Dichlorobenzene	10	10	10	10	10	10	
10	3,3'-Dichlorobenzidine	10	10	10	10	10	10	
10	2,4-Dichlorophenol	10	10	10	10	10	10	
10	Diethylphthalate	10	10	10	10	10	10	
10	2,4-Dimethylphenol	10	10	10	10	10	10	
10	Dimethyl Phthalate	10	10	10	10	10	10	
10	D1-n-butylphthalate	10	10	10	10	10	10	
25	4,6-Dinitro-2-methylphenol	25	25	25	25	25	25	
25	2,4-Dinitrophenol	25	25	25	25	25	25	
10	2,4-Dinitrotoluene	10	10	10	10	10	10	
10	2,6-Dinitrotoluene	10	10	10	10	10	10	
10	D1-n-Octyl Phthalate	10	10	10	10	10	10	
10	bis(2-Ethylhexyl)Phthalate	10	10	10	10	10	10	
10	Fluoranthene	10	10	10	10	10	10	
10	Benzo(b)Fluoranthene	10	10	10	10	10	10	
10	Benzo(k)Fluoranthene	10	10	10	10	10	10	
10	Fluorene	10	10	10	10	10	10	
10	Hexachlorobenzene	10	10	10	10	10	10	
10	Hexachlorobutadiene	10	10	10	10	10	10	
10	Hexachlorocyclopentadiene	10	10	10	10	10	10	
10	Hexachloroethane	10	10	10	10	10	10	

Notes:

All units are ug/l
Sample Quantitation Limit (SQL) is CROL multiplied by the dilution factor and adjusted for percent moisture as appropriate.
The Qual column indicates the qualifier applied to the result following data validation (see below).

Qualifiers:.....
A - Acceptable (Quantitative) Data
J - Estimated (Semi-quantitative) Data
B - Present in associated blanks (Semi-quantitative Data)
K - Biased High (Semi-quantitative) Data
L - Biased Low (Semi-quantitative) Data
UJ - Not Detected/Estimated Data
U - Not Detected/Quantitative Data
UL - Not Detected/Biased Low
N - Tentative Identification
R - Unusable Data
Q - Not Applicable

AR302812

Matrix: Groundwater Well

Semi-volatile Organics

CRQL	Parameter	Sample Point GUM17/18			Sample Point GUM19S/18			Sample Point GUR190/18		
		SQL	Result	Qual	SQL	Result	Qual	SQL	Result	Qual
10	Indeno(1,2,3-cd)Pyrene	10	10	U	10	10	U	10	10	U
10	Isophorone	10	10	U	10	10	U	10	10	U
10	2-Methylnaphthalene	10	10	U	10	10	U	10	10	U
10	2-Methylphenol	10	10	U	10	10	U	10	10	U
10	4-Methylphenol	10	10	U	10	10	U	10	10	U
10	Naphthalene	10	10	U	10	10	U	10	10	U
25	2-Nitroaniline	25	25	U	25	25	U	25	25	U
25	3-Nitroaniline	25	25	U	25	25	U	25	25	U
25	4-Nitroaniline	25	25	U	25	25	U	25	25	U
10	Nitrobenzene	10	10	U	10	10	U	10	10	U
10	2-Nitrophenol	10	10	U	10	10	U	10	10	U
25	4-Nitrophenol	25	25	U	25	25	U	25	25	U
10	N-Nitroso-di-n-propylamine	10	10	U	10	10	U	10	10	U
10	N-Nitrosodiphenylamine	10	10	U	10	10	U	10	10	U
10	2,2-Dybis(1-Chloropropane)	10	10	U	10	10	U	10	10	U
25	Pentachlorophenol	25	25	U	25	25	U	25	25	U
10	Benzo(g,h,i)Perylene	10	10	U	10	10	U	10	10	U
10	Phenanthrene	10	10	U	10	10	U	10	10	U
10	Pyrene	10	10	U	10	10	U	10	10	U
10	Benzo(a)Pyrene	10	10	U	10	10	U	10	10	U
10	1,2,4-Trichlorobenzene	10	10	U	10	10	U	10	10	U
25	2,4,5-Trichlorophenol	25	25	U	25	25	U	25	25	U
10	2,4,6-Trichlorophenol	10	10	U	10	10	U	10	10	U

Notes:
All units are ug/l
Sample Quantitation Limit (SQL) is CRQL multiplied by the dilution factor and adjusted for percent moisture as appropriate.
The Qual column indicates the qualifier applied to the result following data validation (see below).

Qualifiers:
A - Acceptable (Quantitative) Data
J - Estimated (Semi-quantitative) Data
B - Present in associated blanks (Semi-quantitative Data)
K - Biased High (Semi-quantitative) Data
L - Biased Low (Semi-quantitative) Data
UJ - Not Detected/Estimated Data
U - Not Detected/Quantitative Data
UL - Not Detected/Biased Low
N - Tentative Identification
R - Unusable Data
Q - Not Applicable

AR302813

PESTICIDES/PCBs

Phase 1A

AR302814

913
 SUMMARY OF CLP ORGANICS ANALYSES - PHASE 1A
 BERKS LANDFILL REMEDIAL INVESTIGATION
 BERKS COUNTY, PENNSYLVANIA

Matrix: Groundwater Well

Pesticide Organics

CRQL	Parameter	Sample Point GV18S/1A			Sample Point GV18S/ED			Sample Point GVC3D/1A			Sample Point GVC3S/1A		
		SOL	Result	Qual	SOL	Result	Qual	SOL	Result	Qual	SOL	Result	Qual
0.05	Aldrin	0.050	0.050	UL	0.050	0.050	U	0.050	0.050	UL	0.050	0.050	UL
1.0	Aroclor-1016	1.0	1.0	UL	1.0	1.0	U	1.0	1.0	UL	1.0	1.0	UL
2.0	Aroclor-1221	2.0	2.0	UL	2.0	2.0	U	2.0	2.0	UL	2.0	2.0	UL
1.0	Aroclor-1232	1.0	1.0	UL	1.0	1.0	U	1.0	1.0	UL	1.0	1.0	UL
1.0	Aroclor-1242	1.0	1.0	UL	1.0	1.0	U	1.0	1.0	UL	1.0	1.0	UL
1.0	Aroclor-1248	1.0	1.0	UL	1.0	1.0	U	1.0	1.0	UL	1.0	1.0	UL
1.0	Aroclor-1254	1.0	1.0	UL	1.0	1.0	U	1.0	1.0	UL	1.0	1.0	UL
1.0	Aroclor-1260	1.0	1.0	UL	1.0	1.0	U	1.0	1.0	UL	1.0	1.0	UL
0.05	alpha-BHC	0.050	0.050	UL	0.050	0.050	U	0.050	0.050	UL	0.050	0.050	UL
0.05	beta-BHC	0.050	0.050	UL	0.050	0.050	U	0.050	0.050	UL	0.050	0.050	UL
0.05	delta-BHC	0.050	0.050	UL	0.050	0.050	U	0.050	0.050	UL	0.050	0.050	UL
0.05	gamma-BHC (Lindane)	0.050	0.050	UL	0.050	0.050	U	0.050	0.050	UL	0.050	0.050	UL
0.05	alpha-Chlordane	0.050	0.050	UL	0.050	0.050	U	0.050	0.050	UL	0.050	0.050	UL
0.05	gamma-Chlordane	0.050	0.050	UL	0.050	0.050	U	0.050	0.050	UL	0.050	0.050	UL
0.10	4,4'-DDE	0.10	0.10	UL	0.10	0.10	U	0.10	0.10	UL	0.10	0.10	UL
0.10	4,4'-DDD	0.10	0.10	UL	0.10	0.10	U	0.10	0.10	UL	0.10	0.10	UL
0.10	4,4'-DDT	0.10	0.10	UL	0.10	0.10	U	0.10	0.10	UL	0.10	0.10	UL
0.10	Dieldrin	0.10	0.10	UL	0.10	0.10	U	0.10	0.10	UL	0.10	0.10	UL
0.05	Endosulfan I	0.050	0.050	UL	0.050	0.050	U	0.050	0.050	UL	0.050	0.050	UL
0.10	Endosulfan II	0.10	0.10	UL	0.10	0.10	U	0.10	0.10	UL	0.10	0.10	UL
0.10	Endosulfan Sulfate	0.10	0.10	UL	0.10	0.10	U	0.10	0.10	UL	0.10	0.10	UL
0.10	Endrin	0.10	0.10	UL	0.10	0.10	U	0.10	0.10	UL	0.10	0.10	UL
0.10	Endrin Aldehyde	0.10	0.10	UL	0.10	0.10	U	0.10	0.10	UL	0.10	0.10	UL
0.10	Endrin Ketone	0.10	0.10	UL	0.10	0.10	U	0.10	0.10	UL	0.10	0.10	UL
0.05	Heptachlor	0.050	0.050	UL	0.050	0.050	U	0.050	0.050	UL	0.050	0.050	UL
0.05	Heptachlor Epoxide	0.050	0.050	UL	0.050	0.050	U	0.050	0.050	UL	0.050	0.050	UL
0.50	Methoxychlor	0.50	0.50	UL	0.50	0.50	U	0.50	0.50	UL	0.50	0.50	UL
5.0	Toxophene	5.0	5.0	UL	5.0	5.0	U	5.0	5.0	UL	5.0	5.0	UL

Notes:
 All units are ug/l
 Sample Quantitation Limit (SOL) is CRQL multiplied by the dilution factor and adjusted for percent moisture as appropriate.
 The Qual column indicates the qualifier applied to the result following data validation (see below).

Qualifiers:.....
 A - Acceptable (Quantitative) Data
 J - Estimated (Semi-quantitative) Data
 B - Present in associated blanks (Semi-quantitative Data)
 K - Biased High (Semi-quantitative) Data
 L - Biased Low (Semi-quantitative) Data
 UJ - Not Detected/Estimated Data
 U - Not Detected/Quantitative Data
 UL - Not Detected/Biased Low
 N - Tentative Identification
 R - Unusable Data
 Q - Not Applicable

SUMMARY OF CLP ORGANICS ANALYSES - PHASE 1A
 BERKS LANDFILL REMEDIAL INVESTIGATION
 BERKS COUNTY, PENNSYLVANIA

Matrix: Groundwater Well

Pesticide Organics

CROL	Parameter	Sample Point GR180/1A			Sample Point GV-16/1A			Sample Point GV-16/1A			Sample Point GU-C5/1A		
		SQL	Result	Qual	SQL	Result	Qual	SQL	Result	Qual	SQL	Result	Qual
0.05	Aldrin	0.050	0.050	UL	0.050	0.050	UL	0.050	0.050	UL	0.050	0.050	UL
1.0	Aroclor-1016	1.0	1.0	UL	1.0	1.0	UL	1.0	1.0	UL	1.0	1.0	UL
2.0	Aroclor-1221	2.0	2.0	UL	2.0	2.0	UL	2.0	2.0	UL	2.0	2.0	UL
1.0	Aroclor-1232	1.0	1.0	UL	1.0	1.0	UL	1.0	1.0	UL	1.0	1.0	UL
1.0	Aroclor-1242	1.0	1.0	UL	1.0	1.0	UL	1.0	1.0	UL	1.0	1.0	UL
1.0	Aroclor-1248	1.0	1.0	UL	1.0	1.0	UL	1.0	1.0	UL	1.0	1.0	UL
1.0	Aroclor-1254	1.0	1.0	UL	1.0	1.0	UL	1.0	1.0	UL	1.0	1.0	UL
1.0	Aroclor-1260	1.0	1.0	UL	1.0	1.0	UL	1.0	1.0	UL	1.0	1.0	UL
0.05	alpha-BHC	0.050	0.050	UL	0.050	0.050	UL	0.050	0.050	UL	0.050	0.050	UL
0.05	beta-BHC	0.050	0.050	UL	0.050	0.050	UL	0.050	0.050	UL	0.050	0.050	UL
0.05	delta-BHC	0.050	0.050	UL	0.050	0.050	UL	0.050	0.050	UL	0.050	0.050	UL
0.05	gamma-BHC (Lindane)	0.050	0.050	UL	0.050	0.050	UL	0.050	0.050	UL	0.050	0.050	UL
0.05	alpha-Chlordane	0.050	0.050	UL	0.050	0.050	UL	0.050	0.050	UL	0.050	0.050	UL
0.05	gamma-Chlordane	0.050	0.050	UL	0.050	0.050	UL	0.050	0.050	UL	0.050	0.050	UL
0.10	4,4'-DDE	0.10	0.10	UL	0.10	0.10	UL	0.10	0.10	UL	0.10	0.10	UL
0.10	4,4'-DDE	0.10	0.10	UL	0.10	0.10	UL	0.10	0.10	UL	0.10	0.10	UL
0.10	4,4'-DDT	0.10	0.10	UL	0.10	0.10	UL	0.10	0.10	UL	0.10	0.10	UL
0.10	Dieldrin	0.10	0.10	UL	0.10	0.10	UL	0.10	0.10	UL	0.10	0.10	UL
0.05	Endosulfan I	0.050	0.050	UL	0.050	0.050	UL	0.050	0.050	UL	0.050	0.050	UL
0.10	Endosulfan II	0.10	0.10	UL	0.10	0.10	UL	0.10	0.10	UL	0.10	0.10	UL
0.10	Endosulfan Sulfate	0.10	0.10	UL	0.10	0.10	UL	0.10	0.10	UL	0.10	0.10	UL
0.10	Endrin	0.10	0.10	UL	0.10	0.10	UL	0.10	0.10	UL	0.10	0.10	UL
0.10	Endrin Aldehyde	0.10	0.10	UL	0.10	0.10	UL	0.10	0.10	UL	0.10	0.10	UL
0.10	Endrin Ketone	0.10	0.10	UL	0.10	0.10	UL	0.10	0.10	UL	0.10	0.10	UL
0.05	Heptachlor	0.050	0.050	UL	0.050	0.050	UL	0.050	0.050	UL	0.050	0.050	UL
0.05	Heptachlor Epoxide	0.050	0.050	UL	0.050	0.050	UL	0.050	0.050	UL	0.050	0.050	UL
0.50	Methoxychlor	0.50	0.50	UL	0.50	0.50	UL	0.50	0.50	UL	0.50	0.50	UL
5.0	Toxaphene	5.0	5.0	UL	5.0	5.0	UL	5.0	5.0	UL	5.0	5.0	UL

Notes:
 All units are ug/l
 Sample Quantitation Limit (SQL) is CROL multiplied by the dilution factor and adjusted for percent moisture as appropriate.
 The qual column indicates the qualifier applied to the result following data validation (see below).

Qualifiers:
 A - Acceptable (Quantitative) Data
 J - Estimated (Semi-quantitative) Data
 B - Present in associated blanks (Semi-quantitative Data)
 K - Biased High (Semi-quantitative) Data
 L - Biased Low (Semi-quantitative) Data
 UJ - Not Detected/Estimated Data
 U - Not Detected/Quantitative Data
 UL - Not Detected/Biased Low
 N - Tentative Identification
 R - Unusable Data
 Q - Not Applicable

AR302816

THIS PAGE LEFT INTENTIONALLY BLANK

AR302817

SUMMARY OF CLP ORGANIC ANALYSES - PHASE 1A
BERKS LANDFILL REMEDIAL INVESTIGATION
BERKS COUNTY, PENNSYLVANIA

Matrix: Groundwater Well

Pesticide Organics

CROL	Parameter	Sample Point GXG01/1A Date Sampled: 09/02/92 Dilution Factor: 1.0 Percent Moisture: NA %			Sample Point GXG02/1A Date Sampled: 09/02/92 Dilution Factor: 1.0 Percent Moisture: NA %			Sample Point GXG03/1A Date Sampled: 09/02/92 Dilution Factor: 1.0 Percent Moisture: NA %			Sample Point Date Sampled: Dilution Factor: Percent Moisture: %		
		SQL	Result	Qual	SQL	Result	Qual	SQL	Result	Qual	SQL	Result	Qual
0.05	Aldrin	0.050	0.050	UL	0.050	0.050	UL	0.050	0.050	UL			
1.0	Aroclor-1016	1.0	1.0	UL	1.0	1.0	UL	1.0	1.0	UL			
2.0	Aroclor-1221	2.0	2.0	UL	2.0	2.0	UL	2.0	2.0	UL			
1.0	Aroclor-1232	1.0	1.0	UL	1.0	1.0	UL	1.0	1.0	UL			
1.0	Aroclor-1242	1.0	1.0	UL	1.0	1.0	UL	1.0	1.0	UL			
1.0	Aroclor-1248	1.0	1.0	UL	1.0	1.0	UL	1.0	1.0	UL			
1.0	Aroclor-1254	1.0	1.0	UL	1.0	1.0	UL	1.0	1.0	UL			
1.0	Aroclor-1260	1.0	1.0	UL	1.0	1.0	UL	1.0	1.0	UL			
0.05	alpha-BHC	0.050	0.050	UL	0.050	0.050	UL	0.050	0.050	UL			
0.05	beta-BHC	0.050	0.050	UL	0.050	0.050	UL	0.050	0.050	UL			
0.05	delta-BHC	0.050	0.050	UL	0.050	0.050	UL	0.050	0.050	UL			
0.05	gamma-BHC (Lindane)	0.050	0.050	UL	0.050	0.050	UL	0.050	0.050	UL			
0.05	alpha-Chlordane	0.050	0.050	UL	0.050	0.050	UL	0.050	0.050	UL			
0.05	gamma-Chlordane	0.050	0.050	UL	0.050	0.050	UL	0.050	0.050	UL			
0.10	4,4'-DDO	0.10	0.10	UL	0.10	0.10	UL	0.10	0.10	UL			
0.10	4,4'-DDE	0.10	0.10	UL	0.10	0.10	UL	0.10	0.10	UL			
0.10	4,4'-DDT	0.10	0.10	UL	0.10	0.10	UL	0.10	0.10	UL			
0.10	Dieldrin	0.10	0.10	UL	0.10	0.10	UL	0.10	0.10	UL			
0.05	Endosulfan I	0.050	0.050	UL	0.050	0.050	UL	0.050	0.050	UL			
0.10	Endosulfan II	0.10	0.10	UL	0.10	0.10	UL	0.10	0.10	UL			
0.10	Endosulfan Sulfate	0.10	0.10	UL	0.10	0.10	UL	0.10	0.10	UL			
0.10	Endrin	0.10	0.10	UL	0.10	0.10	UL	0.10	0.10	UL			
0.10	Endrin Aldehyde	0.10	0.10	UL	0.10	0.10	UL	0.10	0.10	UL			
0.10	Endrin Ketone	0.10	0.10	UL	0.10	0.10	UL	0.10	0.10	UL			
0.05	Heptachlor	0.050	0.050	UL	0.050	0.050	UL	0.050	0.050	UL			
0.05	Heptachlor Epoxide	0.050	0.050	UL	0.050	0.050	UL	0.050	0.050	UL			
0.50	Methoxychlor	0.50	0.50	UL	0.50	0.50	UL	0.50	0.50	UL			
5.0	Toxaphene	5.0	5.0	UL	5.0	5.0	UL	5.0	5.0	UL			

Notes:
All units are ug/l
Sample Quantitation Limit (SQL) is CROL multiplied by the dilution factor and adjusted for percent moisture as appropriate.
The Qual column indicates the qualifier applied to the result following data validation (see below).

Qualifiers:
A - Acceptable (Quantitative) Data
K - Biased High (Semi-quantitative) Data
J - Estimated (Semi-quantitative) Data
B - Present in associated blanks (Semi-quantitative Data)
L - Biased Low (Semi-quantitative) Data
UL - Not Detected/Estimated Data
UL - Not Detected/Quantitative Data
UL - Not Detected/Biased Low
N - Tentative Identification
R - Unusable Data
Q - Not Applicable

AR302818

TOTAL (UNFILTERED) METALS

Phase 1A
Phase 1B

AR302819

October 1993

913-6773

SUMMARY OF CLP TOTAL INORGANICS ANALYSES - PHASE 1A
BERKS LANDFILL REMEDIAL INVESTIGATION - PHASE 1A
BERKS COUNTY, PENNSYLVANIA

Matrix: Groundwater Well

Parameter	Sample Point GR-18/1A			Sample Point GU-03S/1A			Sample Point GU-05/1A			Sample Point GU-16/1A		
	CRDL	IDL	Result	CRDL	IDL	Result	CRDL	IDL	Result	CRDL	IDL	Result
Aluminum	200	14.0	104	200	14.0	3160	200	14.0	81.8	200	14.0	3330
Antimony	60.0	14.0	14.0	60.0	14.0	16.2	60.0	14.0	14.0	60.0	14.0	14.0
Arsenic	10.0	3.0	4.8	10.0	3.0	6.6	10.0	3.0	3.0	10.0	3.0	20.0
Barium	200	0.40	32.6	200	0.40	116	200	0.40	42.6	200	0.40	389
Beryllium	5.0	0.30	0.30	5.0	0.30	1.0	5.0	0.30	0.30	5.0	0.30	0.30
Cadmium	5.0	1.9	2.4	5.0	1.9	2.2	5.0	1.9	14.4	5.0	1.9	1.9
Calcium	5000	4.1	57100	5000	4.1	251000	5000	4.1	80100	5000	4.1	465000
Chromium	10.0	3.6	3.6	10.0	3.6	3.6	10.0	3.6	3.6	10.0	3.6	13.7
Cobalt	50.0	2.0	2.0	50.0	2.0	15.6	50.0	2.0	4.2	50.0	2.0	10.5
Copper	25.0	1.9	11.6	25.0	1.9	12.6	25.0	1.9	8.0	25.0	1.9	10.8
Iron	100	9.2	146.0	100	9.2	7220	100	9.2	119000	100	9.2	9670
Lead	3.0	1.0	7.4	3.0	1.0	9.6	3.0	1.0	3.4	3.0	1.0	27.3
Magnesium	5000	12.2	9780	5000	12.2	153000	5000	12.2	34400	5000	12.2	129000
Manganese	15.0	0.60	24.5	15.0	0.60	10100	15.0	0.60	1300	15.0	0.60	12200
Mercury	0.20	0.05	0.05	0.20	0.04	0.04	0.20	0.04	0.07	0.20	0.04	0.05
Nickel	40.0	7.9	7.9	40.0	7.9	104	40.0	7.9	10.3	40.0	7.9	32.4
Potassium	5000	64.0	7060	5000	64.0	3720	5000	64.0	5340	5000	64.0	15600
Selenium	5.0	5.0	5.0	5.0	5.0	5.0	5.0	5.0	5.0	5.0	5.0	5.0
Silver	10.0	2.0	2.0	10.0	2.0	2.0	10.0	2.0	5.4	10.0	2.0	2.0
Sodium	5000	8.0	17400	5000	8.0	48200	5000	8.0	23100	5000	8.0	71900
Thallium	10.0	2.0	2.0	10.0	2.0	10.0	10.0	2.0	2.0	10.0	2.0	2.0
Vanadium	50.0	1.8	1.8	50.0	1.8	12.8	50.0	1.8	8.3	50.0	1.8	8.1
Zinc	20.0	1.2	22.0	20.0	1.2	30.9	20.0	1.2	11.0	20.0	1.2	47.4
Cyanide	10.0	10.0	10.0	10.0	10.0	10.0	10.0	10.0	10.0	10.0	10.0	10.0

Notes: All units are ug/l
 Sample Quantitation Limit (SQL) is the Instrument Detection Limit (IDL).
 The CRDLs and IDLs have been adjusted for dilution factors and percent solids as appropriate.

Qualifiers:.....
 A - Acceptable (Quantitative) Data
 J - Estimated (Semi-quantitative) Data
 K - Biased High (Semi-quantitative) Data
 L - Biased Low (Semi-quantitative) Data
 B - Present in associated blanks (Semi-quantitative) Data
 UJ - Not Detected/Estimated Data
 U - Estimated/Quantitative Data
 UL - Not Detected/Biased Low
 M - Tentative Identification
 R - Unusable Data
 Q - Not Applicable

AR302820

**SUMMARY OF CLP TOTAL INORGANICS ANALYSES - PHASE 1A
BERKS LANDFILL REMEDIAL INVESTIGATION - PHASE 1A
BERKS COUNTY, PENNSYLVANIA**

Matrix: Groundwater Well

Parameter	Sample Point GJ-18S/1A			Sample Point GJ-18S/1A			Sample Point GJ-18S/1A			Sample Point GJ-18S/1A			Sample Point GJ-18S/1A			Sample Point GJ-18S/1A				
	CRDL	IDL	Result	Qual	CRDL	IDL	Result	Qual	CRDL	IDL	Result	Qual	CRDL	IDL	Result	Qual	CRDL	IDL	Result	Qual
Aluminum	200	14.0	50.2	B	200	14.0	75.4	B	200	14.0	41.9	B	200	14.0	14.0	J	200	14.0	55700	J
Antimony	60.0	14.0	14.0	U	60.0	14.0	14.0	U	60.0	14.0	14.0	U	60.0	14.0	14.0	U	60.0	14.0	14.0	U
Arsenic	10.0	3.0	3.0	U	10.0	3.0	3.0	U	10.0	3.0	3.0	U	10.0	3.0	3.0	U	10.0	3.0	8.9	A
Barium	200	0.40	58.3	J	200	0.40	60.6	J	200	0.40	56.0	J	200	0.40	309	J	200	0.40	309	J
Beryllium	5.0	0.30	0.30	U	5.0	0.30	0.30	U	5.0	0.30	0.30	U	5.0	0.30	2.4	A	5.0	0.30	2.4	A
Cadmium	5.0	1.9	1.9	U	5.0	1.9	1.9	U	5.0	1.9	1.9	U	5.0	1.9	11.4	A	5.0	1.9	11.4	A
Calcium	5000	4.1	203000	J	5000	4.1	209000	J	5000	4.1	178000	J	5000	4.1	113000	J	5000	4.1	113000	J
Chromium	10.0	3.6	3.6	UL	10.0	3.6	3.6	UL	10.0	3.6	3.6	UL	10.0	3.6	65.5	J	10.0	3.6	65.5	J
Cobalt	50.0	2.0	2.0	U	50.0	2.0	2.0	U	50.0	2.0	2.0	U	50.0	2.0	30.2	A	50.0	2.0	30.2	A
Copper	25.0	1.9	1.9	U	25.0	1.9	4.2	B	25.0	1.9	3.4	B	25.0	1.9	152	A	25.0	1.9	152	A
Iron	100	9.2	82.2	B	100	9.2	114	B	100	9.2	16500	B	100	9.2	52000	J	100	9.2	52000	J
Lead	3.0	1.0	1.6	B	3.0	1.0	2.0	B	3.0	1.0	3.0	B	3.0	1.0	16.4	A	3.0	1.0	16.4	A
Magnesium	5000	12.2	47600	J	5000	12.2	49300	J	5000	12.2	83800	J	5000	12.2	44400	J	5000	12.2	44400	J
Manganese	15.0	0.60	62.3	J	15.0	0.60	65.3	J	15.0	0.60	429	J	15.0	0.60	2960	J	15.0	0.60	2960	J
Mercury	0.20	0.04	0.30	B	0.20	0.04	0.10	B	0.20	0.04	0.04	B	0.20	0.04	0.27	B	0.20	0.04	0.27	B
Nickel	40.0	7.9	7.9	U	40.0	7.9	30.4	U	40.0	7.9	7.9	U	40.0	7.9	60.1	B	40.0	7.9	60.1	B
Potassium	5000	64.0	2540	J	5000	64.0	2690	J	5000	64.0	1890	J	5000	64.0	8690	A	5000	64.0	8690	A
Selenium	5.0	5.0	5.0	U	5.0	5.0	5.0	U	5.0	5.0	5.0	U	5.0	5.0	5.0	U	5.0	5.0	5.0	U
Silver	10.0	2.0	4.4	B	10.0	2.0	3.5	B	10.0	2.0	2.7	B	10.0	2.0	2.6	B	10.0	2.0	2.6	B
Sodium	5000	8.0	15000	J	5000	8.0	15600	J	5000	8.0	14300	J	5000	8.0	12000	J	5000	8.0	12000	J
Thallium	10.0	10.0	10.0	UL	10.0	10.0	2.0	UL	10.0	10.0	2.0	UL	10.0	10.0	2.0	UL	10.0	10.0	2.0	UL
Vanadium	50.0	1.8	1.8	U	50.0	1.8	1.8	U	50.0	1.8	1.8	U	50.0	1.8	111	J	50.0	1.8	111	J
Zinc	20.0	1.2	7.3	B	20.0	1.2	10.4	B	20.0	1.2	7.3	B	20.0	1.2	152	J	20.0	1.2	152	J
Cyanide	10.0	10.0	10.0	U	10.0	10.0	10.0	U	10.0	10.0	10.0	U	10.0	10.0	10.0	U	10.0	10.0	10.0	U

Notes: All units are ug/l
Sample Quantitation Limit (SQL) is the Instrument Detection Limit (IDL).
The CRDLs and IDLs have been adjusted for dilution factors and percent solids as appropriate.

Qualifiers:.....
A - Acceptable (Quantitative) Data
J - Estimated (Semiquantitative) Data
K - Biased High (Semiquantitative) Data
L - Biased Low (Semiquantitative) Data
B - Present in associated blanks (Semiquantitative) Data
UL - Not Detected/Estimated Data
U - Not Detected/Quantitative Data
M - Tentative Identification
N - Unusable Data
O - Not Applicable

AR302821

October 1993

913-6773

SUMMARY OF CLP TOTAL INORGANICS ANALYSES - PHASE 1A
 BERKS LANDFILL REMEDIAL INVESTIGATION - PHASE 1A
 BERKS COUNTY, PENNSYLVANIA

Matrix: Groundwater Well

Parameter	Sample Point GUG-02/1A			Sample Point GUG-03/1A			Sample Point			Sample Point		
	CRDL	IDL	Result	Qual	CRDL	IDL	Result	Qual	CRDL	IDL	Result	Qual
Aluminum	200	14.0	58300	J	200	14.0	44600	J				
Antimony	60.0	14.0	14.0	U	60.0	14.0	14.0	U				
Arsenic	10.0	3.0	8.0	A	10.0	3.0	3.4	L				
Barium	200	0.40	173	J	200	0.40	72.1	J				
Beryllium	5.0	0.30	4.5	A	5.0	0.30	0.30	U				
Cadmium	5.0	1.9	13.1	J	5.0	1.9	9.4	J				
Calcium	5000	4.1	35300	J	5000	4.1	22000	J				
Chromium	10.0	3.6	97.5	J	10.0	3.6	4.4	J				
Cobalt	50.0	2.0	26.7	A	50.0	2.0	26.0	A				
Copper	25.0	1.9	83.7	J	25.0	1.9	272	J				
Iron	100	9.2	66400	J	100	9.2	56400	J				
Lead	3.0	1.0	11.2	J	3.0	1.0	10.0	A				
Magnesium	5000	12.2	22600	J	5000	12.2	9970	J				
Manganese	15.0	0.60	844	J	15.0	0.60	1180	J				
Mercury	0.2	0.04	0.19	B	0.2	0.04	0.21	B				
Nickel	40.0	7.9	75.5	A	40.0	7.9	36.4	B				
Potassium	5000	64.0	3950	J	5000	64.0	2110	J				
Selenium	5.0	5.0	5.0	U	5.0	5.0	5.0	U				
Silver	10.0	2.0	2.0	U	10.0	2.0	2.0	U				
Sodium	5000	8.0	9340	J	5000	8.0	5940	J				
Thallium	10.0	2.0	2.0	U	10.0	2.0	2.0	U				
Vanadium	50.0	1.8	156	J	50.0	1.8	122	J				
Zinc	20.0	1.2	124	J	20.0	1.2	67.4	J				
Cyanide	10.0	10.0	10.0	U	10.0	10.0	10.0	U				

Notes: All units are ug/l

Sample Quantitation Limit (SQL) is the Instrument Detection Limit (IDL).

The CRDLs and IDLs have been adjusted for dilution factors and percent solids as appropriate.

Qualifiers.....
 A - Acceptable (Quantitative) Data
 J - Estimated (Semiquantitative) Data
 K - Biased High (Semiquantitative) Data
 L - Biased Low (Semiquantitative) Data
 B - Present in associated blanks (Semiquantitative) Data
 UJ - Not Detected/Estimated Data
 U - Not Detected/Quantitative Data
 UL - Not Detected/Biased Low
 N - Tentative Identification
 R - Unusable Data
 Q - Not Applicable

AR302822

October 1993

913-6773

SUMMARY OF CLP TOTAL INORGANICS ANALYSES - PHASE 1B
 BERKS LANDFILL REMEDIAL INVESTIGATION - PHASE 1B
 BERKS COUNTY, PENNSYLVANIA

Matrix: Groundwater Well

Parameter	Sample Point GUC01/1B			Sample Point GUC02/1B			Sample Point GUC030/1B			Sample Point GUC030/FD		
	CRDL	IDL	Result	Qual	CRDL	IDL	Result	Qual	CRDL	IDL	Result	Qual
Aluminum	200	44.0	127	B	200	44.0	161	B	200	44.0	660	J
Antimony	60.0	35.0	35.0	UL	60.0	35.0	35.0	UL	60.0	35.0	35.0	UL
Arsenic	10.0	4.0	4.0	UL	10.0	4.0	4.0	UL	10.0	4.0	4.0	UL
Barium	200	2.0	3.1	A	200	2.0	21.2	A	200	2.0	22.7	A
Beryllium	5.0	1.0	1.0	U	5.0	1.0	1.0	U	5.0	1.0	1.0	U
Cadmium	5.0	3.0	3.6	B	5.0	3.0	3.9	B	5.0	3.0	3.0	U
Calcium	5000	17.0	29300	A	5000	17.0	33300	A	5000	17.0	134000	A
Chromium	10.0	8.0	8.0	U	10.0	8.0	8.0	U	10.0	8.0	8.0	U
Chromium	50.0	9.0	9.0	U	50.0	9.0	9.0	U	50.0	9.0	9.0	U
Cobalt	25.0	4.0	4.0	U	25.0	4.0	4.0	U	25.0	4.0	9.7	J
Copper	100	25.0	14200	J	100	25.0	8590	J	100	25.0	862	J
Iron	3.0	2.0	2.0	U	3.0	2.0	2.0	U	3.0	2.0	2.0	U
Lead	5000	48.0	10000	J	5000	48.0	11300	J	5000	48.0	46400	J
Magnesium	15.0	1.0	1020	J	15.0	1.0	111	J	15.0	1.0	145	J
Manganese	0.20	0.20	0.20	U	0.20	0.20	0.20	U	0.20	0.20	0.20	U
Mercury	40.0	18.0	18.0	U	40.0	18.0	18.0	U	40.0	18.0	18.0	U
Nickel	5000	1270	1270	U	5000	1270	1270	U	5000	1270	1870	A
Potassium	5.0	3.0	3.0	R	5.0	3.0	3.0	R	5.0	3.0	3.0	R
Selenium	10.0	4.0	4.0	U	10.0	4.0	4.0	U	10.0	4.0	4.0	U
Silver	5000	123	6990	J	5000	123	9220	J	5000	123	12100	J
Sodium	10.0	3.0	3.0	U	10.0	3.0	3.0	U	10.0	3.0	3.0	U
Thallium	50.0	6.0	6.0	U	50.0	6.0	6.0	U	50.0	6.0	6.0	U
Vanadium	20.0	4.0	16.2	B	20.0	4.0	6.7	B	20.0	4.0	39.1	J
Zinc			Not Analyzed				Not Analyzed				Not Analyzed	
Cyanide			Not Analyzed				Not Analyzed				Not Analyzed	

Notes: All units are ug/l
 Sample Quantitation Limit (SQL) is the Instrument Detection Limit (IDL).
 The CRDLs and IDLs have been adjusted for dilution factors and percent solids as appropriate.

Qualifiers:.....
 A - Acceptable (Quantitative) Data
 J - Estimated (Semi-quantitative) Data
 K - Biased High (Semi-quantitative) Data
 L - Biased Low (Semi-quantitative) Data
 B - Present in associated blanks (Semi-quantitative) Data
 UJ - Not Detected/Estimated Data
 U - Not Detected/Quantitative Data
 UL - Not Detected/Biased Low
 N - Tentative Identification
 R - Unusable Data
 Q - Not Applicable

AR302823

October 1993

913-6773

SUMMARY OF CLP TOTAL INORGANICS ANALYSES - PHASE 1B
 BERKS LANDFILL REMEDIAL INVESTIGATION - PHASE 1B
 BERKS COUNTY, PENNSYLVANIA

Matrix: Groundwater Well

Parameter	Sample Point GMC03S/1B			Sample Point GMC04P/1B			Sample Point GMC04S/1B			Sample Point GMC06P/1B		
	CRDL	IDL	Qual	CRDL	IDL	Qual	CRDL	IDL	Qual	CRDL	IDL	Qual
Aluminum	200	44.0	J	200	44.0	B	200	44.0	B	200	44.0	B
Antimony	60.0	35.0	UL	60.0	35.0	UL	60.0	35.0	UL	60.0	35.0	B
Arsenic	10.0	4.0	J	10.0	4.0	L	10.0	4.0	UL	10.0	4.0	UL
Barium	200	2.0	A	200	2.0	A	200	2.0	A	200	2.0	A
Beryllium	5.0	1.0	UL	5.0	1.0	B	5.0	1.0	U	5.0	1.0	U
Cadmium	5.0	3.0	A	5.0	3.0	U	5.0	3.0	U	5.0	3.0	U
Calcium	5000	17.0	A	5000	17.0	A	5000	17.0	A	5000	17.0	A
Chromium	10.0	8.0	J	10.0	8.0	U	10.0	8.0	U	10.0	8.0	U
Chromium	10.0	8.0	J	10.0	8.0	U	10.0	8.0	U	10.0	8.0	U
Cobalt	50.0	9.0	A	50.0	9.0	A	50.0	9.0	U	50.0	9.0	U
Copper	25.0	4.0	J	25.0	4.0	J	25.0	4.0	J	25.0	4.0	J
Iron	100	25.0	J	100	25.0	J	100	25.0	J	100	25.0	J
Lead	3.0	2.0	U	3.0	2.0	UL	3.0	2.0	UL	3.0	2.0	UL
Magnesium	5000	48.0	J	5000	48.0	J	5000	48.0	J	5000	48.0	J
Manganese	15.0	1.0	J	15.0	1.0	J	15.0	1.0	J	15.0	1.0	J
Mercury	0.20	0.20	U	0.20	0.20	U	0.20	0.20	U	0.20	0.20	U
Nickel	40.0	18.0	L	40.0	18.0	U	40.0	18.0	U	40.0	18.0	U
Potassium	5000	1270	A	5000	1270	A	5000	1270	A	5000	1270	U
Selenium	5.0	3.0	R	5.0	3.0	R	5.0	3.0	R	5.0	3.0	R
Silver	10.0	4.0	U	10.0	4.0	U	10.0	4.0	U	10.0	4.0	U
Sodium	5000	123	J	5000	123	J	5000	123	J	5000	123	J
Thallium	10.0	3.0	UL	10.0	3.0	UL	10.0	3.0	UL	10.0	3.0	UL
Vanadium	50.0	6.0	U	50.0	6.0	U	50.0	6.0	U	50.0	6.0	U
Zinc	20.0	4.0	J	20.0	4.0	J	20.0	4.0	J	20.0	4.0	J
Cyanide			Not Analyzed			Not Analyzed			Not Analyzed			Not Analyzed
			Result			Result			Result			Result
			1930			506			201			184
			35.0			35.0			35.0			42.9
			5.6			13.6			4.0			4.0
			89.4			79.4			403			52.0
			1.0			3.5			1.0			1.0
			3.3			3.0			3.4			3.0
			210000			37700			145000			35800
			19.8			8.0			8.0			8.0
			14.3			12.7			9.0			9.0
			7.6			10.0			9.5			6.2
			4920			371			2060			133
			2.0			2.0			2.0			2.0
			135000			7260			46100			11300
			7580			361			942			15.7
			0.20			0.20			0.20			0.20
			56.7			18.0			18.0			18.0
			3180			3350			5710			1270
			3.0			15.0			15.0			3.0
			4.0			4.0			4.0			4.0
			36100			35000			18200			7830
			3.0			3.0			3.0			3.0
			6.0			6.0			6.0			10.1
			27.1			81.1			83.9			16.9
			Not Analyzed			Not Analyzed			Not Analyzed			Not Analyzed

Notes: All units are ug/l
 Sample Quantitation Limit (SQL) is the Instrument Detection Limit (IDL).
 The CRDLs and IDLs have been adjusted for dilution factors and percent solids as appropriate.

Qualifiers.....
 A - Acceptable (Quantitative) Data
 J - Estimated (Semi-quantitative) Data
 K - Biased High (Semi-quantitative) Data
 L - Biased Low (Semi-quantitative) Data
 B - Present in associated blanks (Semi-quantitative) Data
 UJ - Not Detected/Estimated Data
 U - Not Detected/Quantitative Data
 UL - Not Detected/Biased Low
 N - Tentative Identification
 R - Unusable Data
 Q - Not Applicable

AR302824

October 1993

913-6773

SUMMARY OF CLP TOTAL INORGANICS ANALYSES - PHASE 1B
 BERKS LANDFILL REMEDIAL INVESTIGATION - PHASE 1B
 BERKS COUNTY, PENNSYLVANIA

Matrix: Groundwater Well

Parameter	Sample Point GUC06S/1B			Sample Point GUC07D/1B			Sample Point GUC07S/1B			Sample Point GUC02/1B		
	CRDL	IDL	Result	CRDL	IDL	Result	CRDL	IDL	Result	CRDL	IDL	Result
Aluminum	200	44.0	230	200	44.0	218	200	64.0	413	200	44.0	7050
Antimony	60.0	35.0	50.7	60.0	35.0	35.0	60.0	35.0	35.1	60.0	35.0	35.0
Arsenic	10.0	4.0	4.0	10.0	4.0	4.2	10.0	4.0	4.0	10.0	4.0	4.4
Barium	200	2.0	107	200	2.0	59.6	200	2.0	323	200	2.0	449
Beryllium	5.0	1.0	1.0	5.0	1.0	1.0	5.0	1.0	1.0	5.0	1.0	1.0
Cadmium	3.0	3.0	3.0	3.0	3.0	5.3	3.0	3.0	3.8	3.0	3.0	3.0
Calcium	5000	17.0	26600	5000	17.0	31500	5000	17.0	333000	5000	17.0	123000
Chromium	10.0	8.0	8.0	10.0	8.0	9.0	10.0	8.0	8.0	10.0	8.0	8.9
Cobalt	50.0	9.0	9.0	50.0	9.0	9.0	50.0	9.0	9.0	50.0	9.0	15.5
Copper	25.0	4.0	4.0	25.0	4.0	4.0	25.0	4.0	9.4	25.0	4.0	31.4
Iron	100	25.0	2980	100	25.0	126	100	25.0	518	100	25.0	32800
Lead	3.0	2.0	2.0	3.0	2.0	2.0	3.0	2.0	2.0	3.0	2.0	12.0
Magnesium	5000	48.0	7200	5000	48.0	6300	5000	48.0	68200	5000	48.0	47100
Manganese	15.0	1.0	64.6	15.0	1.0	42.2	15.0	1.0	1400	15.0	1.0	2510
Mercury	0.20	0.20	0.20	0.20	0.20	0.20	0.20	0.20	0.20	0.20	0.20	0.20
Nickel	40.0	18.0	18.0	40.0	18.0	18.0	40.0	18.0	19.1	40.0	18.0	18.0
Potassium	5000	1270	1640	5000	1270	1850	5000	1270	4060	5000	1270	7540
Selenium	5.0	3.0	3.0	5.0	3.0	3.0	5.0	3.0	3.0	5.0	3.0	30.0
Silver	10.0	4.0	4.0	10.0	4.0	4.0	10.0	4.0	4.0	10.0	4.0	4.0
Sodium	5000	123	7210	5000	123	9650	5000	123	24200	5000	123	33900
Thallium	10.0	3.0	3.0	10.0	3.0	3.0	10.0	3.0	3.0	10.0	3.0	3.0
Vanadium	50.0	6.0	6.0	50.0	6.0	6.0	50.0	6.0	6.0	50.0	6.0	13.8
Zinc	20.0	4.0	11.6	20.0	4.0	7.7	20.0	4.0	29.6	20.0	4.0	70.7
Cyanide			Not Analyzed			Not Analyzed			Not Analyzed			Not Analyzed

Notes: All units are ug/l
 Sample Quantitation Limit (SQL) is the Instrument Detection Limit (IDL).
 The CRDLs and IDLs have been adjusted for dilution factors and percent solids as appropriate.

Qualifiers.....
 A - Acceptable (Quantitative) Data
 J - Estimated (Semi-quantitative) Data
 K - Biased High (Semi-quantitative) Data
 L - Biased Low (Semi-quantitative) Data
 B - Present in associated blanks (Semi-quantitative) Data
 UJ - Not Detected/Estimated Data
 U - Not Detected/Quantitative Data
 UL - Not Detected/Biased Low
 M - Tentative Identification
 R - Unusable Data
 Q - Not Applicable

AR302825

October 1993

913-6773

SUMMARY OF CLP TOTAL INORGANICS ANALYSES - PHASE 1B
 BERKS LANDFILL REMEDIAL INVESTIGATION - PHASE 1B
 BERKS COUNTY, PENNSYLVANIA

Matrix: Groundwater Well

Parameter	Sample Point GMG01/1B			Sample Point GMG02/1B			Sample Point GMG03/1B			Sample Point GMG04/1B		
	CRDL	IDL	Result	CRDL	IDL	Result	CRDL	IDL	Result	CRDL	IDL	Result
Aluminum	200	44.0	13400	200	44.0	29200	200	44.0	5150	200	44.0	362
Antimony	60.0	35.0	35.0	60.0	35.0	35.0	60.0	35.0	35.0	60.0	35.0	46.7
Arsenic	10.0	4.0	5.8	10.0	4.0	7.9	10.0	4.0	4.0	10.0	4.0	21.2
Barium	200	2.0	154	200	2.0	116	200	2.0	11.3	200	2.0	63.0
Beryllium	5.0	1.0	1.0	5.0	1.0	2.2	5.0	1.0	1.0	5.0	1.0	1.0
Cadmium	5.0	3.0	3.0	5.0	3.0	3.0	5.0	3.0	3.0	5.0	3.0	3.0
Calcium	5000	17.0	22400	5000	17.0	37100	5000	17.0	20200	5000	17.0	246000
Chromium	10.0	8.0	15.7	10.0	8.0	57.0	10.0	8.0	8.0	10.0	8.0	8.0
Chromium	50.0	9.0	10.3	50.0	9.0	25.2	50.0	9.0	9.0	50.0	9.0	9.0
Cobalt	25.0	4.0	46.5	25.0	4.0	57.3	25.0	4.0	41.1	25.0	4.0	12.6
Copper	100	25.0	11000	100	25.0	35600	100	25.0	6460	100	25.0	1220
Iron	3.0	2.0	5.1	3.0	2.0	11.9	3.0	2.0	2.5	3.0	2.0	2.2
Lead	5000	48.0	14600	5000	48.0	16600	5000	48.0	6940	5000	48.0	41000
Magnesium	15.0	1.0	6670	15.0	1.0	813	15.0	1.0	158	15.0	1.0	1580
Manganese	0.20	0.20	0.20	0.20	0.20	0.20	0.20	0.20	0.20	0.20	0.20	0.20
Mercury	40.0	18.0	18.0	40.0	18.0	18.0	40.0	18.0	18.0	40.0	18.0	18.0
Nickel	5000	1270	2840	5000	1270	2590	5000	1270	1270	5000	1270	4020
Potassium	5.0	3.0	3.0	5.0	3.0	30.0	5.0	3.0	30.0	5.0	3.0	3.0
Selenium	10.0	4.0	4.0	10.0	4.0	4.0	10.0	4.0	4.0	10.0	4.0	4.0
Silver	5000	123	12200	5000	123	10300	5000	123	5410	5000	123	38200
Sodium	50.0	3.0	3.0	50.0	3.0	3.0	50.0	3.0	3.0	50.0	3.0	3.0
Thallium	50.0	6.0	25.4	50.0	6.0	84.2	50.0	6.0	16.3	50.0	6.0	6.0
Vanadium	20.0	4.0	44.3	20.0	4.0	136	20.0	4.0	25.4	20.0	4.0	68.3
Zinc			Not Analyzed			Not Analyzed			Not Analyzed			Not Analyzed
Cyanide			Not Analyzed			Not Analyzed			Not Analyzed			Not Analyzed

Notes: All units are ug/l
 Sample Quantitation Limit (SQL) is the Instrument Detection Limit (IDL).
 The CRDLs and IDLs have been adjusted for dilution factors and percent solids as appropriate.

Qualifiers.....
 A - Acceptable (Quantitative) Data
 J - Estimated (Semi-quantitative) Data
 K - Biased High (Semi-quantitative) Data
 L - Biased Low (Semi-quantitative) Data
 B - Present in associated blanks (Semi-quantitative) Data
 UJ - Not Detected/Estimated Data
 U - Not Detected/Quantitative Data
 UL - Not Detected/Biased Low
 N - Tentative Identification
 R - Unusable Data
 Q - Not Applicable

AR302826

SUMMARY OF CLP TOTAL INORGANICS ANALYSES - PHASE 1B
 BERKS LANDFILL REMEDIAL INVESTIGATION - PHASE 1B
 BERKS COUNTY, PENNSYLVANIA

Matrix: Groundwater Well

Parameter	Sample Point GMG05/1B			Sample Point GMG06/1B			Sample Point GMG07/1B			Sample Point GMG08/1B		
	CRDL	IDL	Result	CRDL	IDL	Result	CRDL	IDL	Result	CRDL	IDL	Result
Aluminum	200	44.0	18300	200	44.0	186	200	44.0	7510	200	44.0	14100
Antimony	60.0	35.0	35.0	60.0	35.0	35.0	60.0	35.0	35.0	60.0	35.0	35.0
Arsenic	10.0	4.0	10.4	10.0	4.0	4.0	10.0	4.0	7.4	10.0	4.0	4.0
Barium	200	2.0	605	200	2.0	2.0	200	2.0	28.0	200	2.0	73.6
Beryllium	5.0	1.0	1.4	5.0	1.0	1.0	5.0	1.0	1.0	5.0	1.0	1.0
Cadmium	5.0	3.0	3.0	5.0	3.0	3.0	5.0	3.0	3.0	5.0	3.0	3.0
Calcium	5000	17.0	229000	5000	17.0	35500	5000	17.0	26400	5000	17.0	29600
Chromium	10.0	8.0	16.6	10.0	8.0	8.0	10.0	8.0	8.0	10.0	8.0	8.0
Chromium	50.0	9.0	9.0	50.0	9.0	9.0	50.0	9.0	13.8	50.0	9.0	20.1
Cobalt	25.0	4.0	7.0	25.0	4.0	10.8	25.0	4.0	47.6	25.0	4.0	142
Copper	100	25.0	7510	100	25.0	339	100	25.0	16000	100	25.0	30000
Iron	3.0	2.0	12.6	3.0	2.0	2.0	3.0	2.0	4.0	3.0	2.0	7.0
Lead	5000	48.0	48900	5000	48.0	13100	5000	48.0	4290	5000	48.0	7490
Magnesium	15.0	1.0	1260	15.0	1.0	23.3	15.0	1.0	126	15.0	1.0	611
Mercury	0.20	0.20	0.20	0.20	0.20	0.20	0.20	0.20	0.20	0.20	0.20	0.20
Nickel	40.0	18.0	18.0	40.0	18.0	18.0	40.0	18.0	18.0	40.0	18.0	18.0
Potassium	5000	1270	7320	5000	1270	1270	5000	1270	1850	5000	1270	2820
Selenium	50.0	30.0	30.0	50.0	30.0	30.0	50.0	30.0	30.0	50.0	30.0	30.0
Silver	10.0	4.0	4.0	10.0	4.0	4.0	10.0	4.0	4.0	10.0	4.0	4.0
Sodium	5000	123	61500	5000	123	6810	5000	123	12300	5000	123	6780
Thallium	100	30.0	30.0	100	30.0	3.0	100	30.0	3.0	100	30.0	3.0
Vanadium	50.0	6.0	6.0	50.0	6.0	8.5	50.0	6.0	134	50.0	6.0	49.6
Zinc	20.0	4.0	48.6	20.0	4.0	23.8	20.0	4.0	37.8	20.0	4.0	63.5
Cyanide			Not Analyzed			Not Analyzed			Not Analyzed			Not Analyzed

Notes: All units are ug/l
 Sample Quantitation Limit (SQL) is the Instrument Detection Limit (IDL).
 The CRDLs and IDLs have been adjusted for dilution factors and percent solids as appropriate.

Qualifiers.....
 A - Acceptable (Quantitative) Data
 J - Estimated (Semi-quantitative) Data
 K - Biased High (Semi-quantitative) Data
 L - Biased Low (Semi-quantitative) Data
 B - Present in associated blanks (Semi-quantitative) Data
 UL - Not Detected/Estimated Data
 U - Not Detected/Quantitative Data
 UL - Not Detected/Biased Low
 M - Tentative Identification
 R - Unusable Data
 Q - Not Applicable

October 1993

913-6773

SUMMARY OF CLP TOTAL INORGANICS ANALYSES - PHASE 1B
 BERKS LANDFILL REMEDIAL INVESTIGATION - PHASE 1B
 BERKS COUNTY, PENNSYLVANIA

Matrix: Groundwater Well

Parameter	Sample Point GMG10/1B			Sample Point GMG11/1B			Sample Point GMG11/FD			Sample Point GMG12/1B		
	CRDL	IDL	Result	Qual	CRDL	IDL	Result	Qual	CRDL	IDL	Result	Qual
Aluminum	200	44.0	47300	J	200	44.0	520	B	200	44.0	650	J
Antimony	60.0	35.0	35.0	UL	60.0	35.0	46.8	B	60.0	35.0	50.8	UL
Arsenic	10.0	4.0	7.9	J	10.0	4.0	4.0	UL	10.0	4.0	5.3	L
Barium	200	2.0	144	A	200	2.0	1360	A	200	2.0	1380	A
Beryllium	5.0	1.0	5.7	L	5.0	1.0	1.0	U	5.0	1.0	1.0	U
Cadmium	5.0	3.0	3.0	U	5.0	3.0	4.2	B	5.0	3.0	3.2	U
Calcium	5000	17.0	32000	A	5000	17.0	473000	A	5000	17.0	480000	A
Chromium	10.0	8.0	57.5	J	10.0	8.0	8.0	U	10.0	8.0	8.0	U
Cobalt	50.0	9.0	27.5	A	50.0	9.0	9.0	U	50.0	9.0	9.0	U
Copper	25.0	4.0	175	J	25.0	4.0	4.6	J	25.0	4.0	10.8	J
Iron	100	25.0	44000	J	100	25.0	917	J	100	25.0	994	J
Lead	3.0	2.0	17.3	J	3.0	2.0	2.0	UL	3.0	2.0	2.9	J
Magnesium	5000	48.0	22600	J	5000	48.0	107000	J	5000	48.0	109000	J
Manganese	15.0	1.0	1240	J	15.0	1.0	5170	J	15.0	1.0	5250	J
Mercury	0.20	0.20	0.20	U	0.20	0.20	0.20	U	0.20	0.20	0.20	U
Nickel	40.0	18.0	23.9	L	40.0	18.0	51.4	A	40.0	18.0	52.1	A
Potassium	5000	1270	4090	A	5000	1270	8250	A	5000	1270	8830	A
Selenium	50.0	30.0	30.0	R	5.0	3.0	3.0	R	50.0	30.0	3.0	R
Silver	10.0	4.0	4.0	U	10.0	4.0	4.0	U	10.0	4.0	4.0	U
Sodium	5000	123	10000	J	5000	123	151000	J	5000	123	154000	J
Thallium	10.0	3.0	3.0	UL	100	30.0	30.0	UL	100	30.0	30.0	UL
Vanadium	50.0	6.0	107	A	50.0	6.0	6.0	U	50.0	6.0	6.0	U
Zinc	20.0	4.0	116	J	20.0	4.0	18.6	B	20.0	4.0	18.6	B
Cyanide			Not Analyzed				Not Analyzed				Not Analyzed	

Notes: All units are ug/l
 Sample Quantitation Limit (SQL) is the Instrument Detection Limit (IDL).
 The CRDLs and IDLs have been adjusted for dilution factors and percent solids as appropriate.

Qualifiers.....
 A - Acceptable (Quantitative) Data
 J - Estimated (Semi-quantitative) Data
 K - Biased High (Semi-quantitative) Data
 L - Biased Low (Semi-quantitative) Data
 B - Present in associated blends (Semi-quantitative) Data
 UL - Not Detected/Estimated Data
 U - Not Detected/Quantitative Data
 UL - Not Detected/Biased Low
 N - Tentative Identification
 R - Unusable Data
 Q - Not Applicable

AR302828

October 1993

913-6773

SUMMARY OF CLP TOTAL INORGANICS ANALYSES - PHASE 1B
 BERKS LANDFILL REMEDIAL INVESTIGATION - PHASE 1B
 BERKS COUNTY, PENNSYLVANIA

Matrix: Groundwater Well

Parameter	Sample Point GMG12/FD			Sample Point GMG13/1B			Sample Point GMG03/1B			Sample Point GMG06/1B		
	CRDL	IDL	Result	Qual	CRDL	IDL	Result	Qual	CRDL	IDL	Result	Qual
Aluminum	200	44.0	48100	J	200	44.0	136	B	200	44.0	2240	J
Antimony	60.0	35.0	35.0	UL	60.0	35.0	35.0	UL	60.0	35.0	35.0	UL
Arsenic	10.0	4.0	11.3	J	10.0	4.0	16.4	L	10.0	4.0	12.6	J
Barium	200	2.0	615	A	200	2.0	10.0	A	200	2.0	127	A
Beryllium	5.0	1.0	1.9	L	5.0	1.0	1.0	U	5.0	1.0	1.0	U
Cadmium	5.0	3.0	3.0	U	5.0	3.0	3.4	B	5.0	3.0	3.0	U
Calcium	5000	17.0	517000	A	5000	17.0	47400	A	5000	17.0	176000	A
Chromium	10.0	8.0	54.0	J	10.0	8.0	8.0	U	10.0	8.0	8.0	U
Chromium	50.0	9.0	30.4	A	50.0	9.0	9.0	U	50.0	9.0	9.0	A
Cobalt	25.0	4.0	93.0	J	25.0	4.0	4.0	U	25.0	4.0	14.6	J
Copper	100	25.0	42800	J	100	25.0	280	J	100	25.0	10800	J
Iron	30.0	20.0	93.4	J	3.0	2.0	2.0	UL	3.0	2.0	5.9	J
Lead	5000	48.0	190000	J	5000	48.0	13200	J	5000	48.0	101000	J
Magnesium	15.0	1.0	3380	J	15.0	1.0	26.6	J	15.0	1.0	13500	J
Manganese	0.20	0.20	0.20	U	0.20	0.20	0.20	U	0.20	0.20	0.20	U
Mercury	40.0	18.0	38.3	L	40.0	18.0	18.0	U	40.0	18.0	57.0	L
Nickel	5000	1270	11000	A	5000	1270	1270	U	5000	1270	3590	A
Potassium	5.0	3.0	30.0	R	5.0	3.0	3.0	R	5.0	3.0	3.0	R
Selenium	10.0	4.0	4.0	U	10.0	4.0	4.0	U	10.0	4.0	4.0	U
Silver	5000	123	26900	J	5000	123	12100	J	5000	123	188000	J
Sodium	100	30.0	30.0	UL	10.0	3.0	3.0	UL	10.0	3.0	3.0	UL
Thallium	50.0	6.0	54.8	A	50.0	6.0	9.8	A	50.0	6.0	6.0	U
Vanadium	20.0	4.0	236	J	20.0	4.0	10.2	B	20.0	4.0	315	J
Zinc			Not Analyzed				Not Analyzed				Not Analyzed	
Cyanide			Not Analyzed				Not Analyzed				Not Analyzed	
			147000	J			147000	J			147000	J
			35.0	UL			35.0	UL			35.0	UL
			28.1	J			28.1	J			28.1	J
			493	A			493	A			493	A
			11.6	L			11.6	L			11.6	L
			4.1	A			4.1	A			4.1	A
			87100	A			87100	A			87100	A
			8.0	U			8.0	U			8.0	U
			13.8	A			13.8	A			13.8	A
			24.3	J			24.3	J			24.3	J
			75900	J			75900	J			75900	J
			233	J			233	J			233	J
			50300	J			50300	J			50300	J
			582	J			582	J			582	J
			0.20	U			0.20	U			0.20	U
			18.0	UL			18.0	UL			18.0	UL
			4020	A			4020	A			4020	A
			3.0	R			3.0	R			3.0	R
			4.0	U			4.0	U			4.0	U
			105000	J			105000	J			105000	J
			3.0	UL			3.0	UL			3.0	UL
			16.1	A			16.1	A			16.1	A
			919	J			919	J			919	J
			Not Analyzed				Not Analyzed				Not Analyzed	

Notes: All units are ug/l
 Sample quantitation limit (SQL) is the Instrument Detection Limit (IDL).
 The CRDLs and IDLs have been adjusted for dilution factors and percent solids as appropriate.

Qualifiers.....
 A - Acceptable (Quantitative) Data
 J - Estimated (Semi-quantitative) Data
 K - Biased High (Semi-quantitative) Data
 L - Biased Low (Semi-quantitative) Data
 B - Present in associated blanks (Semi-quantitative) Data
 UJ - Not Detected/Estimated Data
 U - Not Detected/Quantitative Data
 UL - Not Detected/Biased Low
 M - Tentative Identification
 R - Unusable Data
 Q - Not Applicable

AR302829

October 1993

913-6773

SUMMARY OF CLP TOTAL INORGANICS ANALYSES - PHASE 1B
 BERKS LANDFILL REMEDIAL INVESTIGATION - PHASE 1B
 BERKS COUNTY, PENNSYLVANIA

Matrix: Groundwater Well

Parameter	Sample Point GM11/1B Date Sampled: 06/30/93 Dilution Factor: NA % Percent Solids: NA %			Sample Point GM140/1B Date Sampled: 07/12/93 Dilution Factor: NA % Percent Solids: NA %			Sample Point GM14S/1B Date Sampled: 07/09/93 Dilution Factor: NA % Percent Solids: NA %			Sample Point GM15D/1B Date Sampled: 07/08/93 Dilution Factor: NA % Percent Solids: NA %		
	CRDL	IDL	Qual	CRDL	IDL	Qual	CRDL	IDL	Qual	CRDL	IDL	Qual
Aluminum	200	44.0	J	200	44.0	B	200	44.0	B	200	44.0	B
Antimony	60.0	35.0	UL	60.0	35.0	UL	60.0	35.0	B	60.0	35.0	UL
Arsenic	10.0	4.0	UL	10.0	4.0	UL	10.0	4.0	L	10.0	4.0	UL
Barium	200	2.0	A	200	2.0	A	200	2.0	A	200	2.0	A
Beryllium	5.0	1.0	U	5.0	1.0	U	5.0	1.0	A	5.0	1.0	A
Cadmium	3.0	3.0	U	3.0	3.0	U	3.0	3.0	B	3.0	3.0	B
Calcium	5000	17.0	A	5000	17.0	A	5000	17.0	A	5000	17.0	A
Chromium	10.0	8.0	J	10.0	8.0	U	10.0	8.0	A	10.0	8.0	U
Cobalt	50.0	9.0	J	50.0	9.0	U	50.0	9.0	U	50.0	9.0	U
Copper	25.0	4.0	J	25.0	4.0	U	25.0	4.0	U	25.0	4.0	J
Iron	100	25.0	J	100	25.0	J	100	25.0	J	100	25.0	J
Lead	3.0	2.0	UL	3.0	2.0	J	3.0	2.0	U	3.0	2.0	J
Magnesium	5000	48.0	U	5000	48.0	J	5000	48.0	J	5000	48.0	J
Manganese	15.0	1.0	J	15.0	1.0	J	15.0	1.0	J	15.0	1.0	J
Mercury	0.20	0.20	U	0.20	0.20	U	0.20	0.20	U	0.20	0.20	U
Nickel	40.0	18.0	U	40.0	18.0	A	40.0	18.0	A	40.0	18.0	U
Potassium	5000	1270	A	5000	1270	A	5000	1270	A	5000	1270	A
Selenium	5.0	3.0	R	5.0	3.0	R	5.0	3.0	R	5.0	3.0	R
Silver	10.0	4.0	U	10.0	4.0	U	10.0	4.0	U	10.0	4.0	U
Sodium	5000	123	J	5000	123	J	5000	123	J	5000	123	J
Thallium	10.0	3.0	UL	10.0	3.0	UL	10.0	3.0	UL	10.0	3.0	UL
Vanadium	50.0	6.0	U	50.0	6.0	U	50.0	6.0	U	50.0	6.0	U
Zinc	20.0	4.0	B	20.0	4.0	J	20.0	4.0	B	20.0	4.0	B
Cyanide			Not Analyzed			Not Analyzed			Not Analyzed			Not Analyzed

Notes: All units are ug/l
 Sample Quantitation Limit (SQL) is the Instrument Detection Limit (IDL).
 The CRDLs and IDLs have been adjusted for dilution factors and percent solids as appropriate.

Qualifiers.....
 A - Acceptable (Quantitative) Data
 J - Estimated (Semi-quantitative) Data
 K - Biased High (Semi-quantitative) Data
 L - Biased Low (Semi-quantitative) Data
 B - Present in associated blanks (Semi-quantitative) Data
 UJ - Not Detected/Estimated Data
 U - Not Detected/Quantitative Data
 UL - Not Detected/Biased Low
 M - Tentative Identification
 R - Unusable Data
 Q - Not Applicable

AR302830

October 1993

913-6773

SUMMARY OF CLP TOTAL INORGANICS ANALYSES - PHASE 1B
 BERKS LANDFILL REMEDIAL INVESTIGATION - PHASE 1B
 BERKS COUNTY, PENNSYLVANIA

Matrix: Groundwater Well

Parameter	Sample Point GUM155/1B			Sample Point GUM16/1B			Sample Point GUM17/1B			Sample Point GUM195/1B		
	CRDL	IDL	Result	CRDL	IDL	Result	CRDL	IDL	Result	CRDL	IDL	Result
Aluminum	200	44.0	18100	200	44.0	873	200	44.0	464	200	44.0	647
Antimony	60.0	35.0	35.0	60.0	35.0	35.0	60.0	35.0	35.0	60.0	35.0	35.0
Arsenic	10.0	4.0	14.7	10.0	4.0	4.0	10.0	4.0	4.0	10.0	4.0	4.0
Barium	200	2.0	330	200	2.0	365	200	2.0	7.7	200	2.0	61.8
Beryllium	5.0	1.0	1.4	5.0	1.0	1.0	5.0	1.0	1.0	5.0	1.0	1.0
Cadmium	5.0	3.0	3.0	5.0	3.0	3.0	5.0	3.0	3.0	5.0	3.0	3.0
Calcium	5000	17.0	88200	5000	17.0	394000	5000	17.0	13400	5000	17.0	384000
Chromium	10.0	8.0	51.5	10.0	8.0	8.0	10.0	8.0	8.0	10.0	8.0	8.0
Chromium	50.0	9.0	29.2	50.0	9.0	9.7	50.0	9.0	9.0	50.0	9.0	9.0
Cobalt	25.0	4.0	24.9	25.0	4.0	4.9	25.0	4.0	5.9	25.0	4.0	6.6
Copper	100	25.0	24500	100	25.0	4110	100	25.0	5400	100	25.0	636
Iron	3.0	2.0	12.7	3.0	2.0	3.8	3.0	2.0	2.1	3.0	2.0	2.0
Lead	5000	48.0	25800	5000	48.0	102000	5000	48.0	5460	5000	48.0	46400
Magnesium	15.0	1.0	27600	15.0	1.0	12600	15.0	1.0	325	15.0	1.0	693
Manganese	0.20	0.20	0.20	0.20	0.20	0.20	0.20	0.20	0.20	0.20	0.20	0.20
Mercury	40.0	18.0	30.2	40.0	18.0	24.6	40.0	18.0	18.0	40.0	18.0	18.0
Nickel	5000	1270	4880	5000	1270	13000	5000	1270	1270	5000	1270	4450
Potassium	50.0	30.0	30.0	50.0	30.0	3.0	50.0	30.0	3.0	50.0	30.0	3.9
Selenium	10.0	4.0	4.0	10.0	4.0	4.0	10.0	4.0	4.0	10.0	4.0	4.0
Silver	5000	123	28600	5000	123	78100	5000	123	3450	5000	123	9570
Sodium	10.0	3.0	3.0	10.0	3.0	30.0	10.0	3.0	3.0	10.0	3.0	3.0
Thallium	50.0	6.0	47.2	50.0	6.0	6.0	50.0	6.0	6.0	50.0	6.0	10.0
Vanadium	20.0	4.0	102	20.0	4.0	16.0	20.0	4.0	47.9	20.0	4.0	17.2
Zinc			Not Analyzed			Not Analyzed			Not Analyzed			Not Analyzed
Cyanide			Not Analyzed			Not Analyzed			Not Analyzed			Not Analyzed

Notes: All units are ug/l
 Sample Quantitation Limit (SQL) is the Instrument Detection Limit (IDL).
 The CRDLs and IDLs have been adjusted for dilution factors and percent solids as appropriate.

Qualifiers.....
 A - Acceptable (Quantitative) Data
 J - Estimated (Semi-quantitative) Data
 K - Biased High (Semi-quantitative) Data
 L - Biased Low (Semi-quantitative) Data
 B - Present in associated blanks (Semi-quantitative) Data
 UJ - Not Detected/Estimated Data
 U - Not Detected/Quantitative Data
 UL - Not Detected/Biased Low
 M - Tentative Identification
 R - Unusable Data
 Q - Not Applicable

AR302831

stot1b.rpt
 4/93|12:41:07

Go Associates

Page

October 1993

913-6773

**SUMMARY OF CLP TOTAL INORGANICS ANALYSES - PHASE 1B
BERKS LANDFILL REMEDIAL INVESTIGATION - PHASE 1B
BERKS COUNTY, PENNSYLVANIA**

Matrix: Groundwater Well

Parameter	Sample Point GWR190/1B		Sample Point		Sample Point		Sample Point		Sample Point			
	CRDL	IDL	Result	Qual	CRDL	IDL	Result	Qual	CRDL	IDL	Result	Qual
Aluminum	200	44.0	40900	J								
Antimony	60.0	35.0	35.0	UL								
Arsenic	10.0	4.0	11.4	J								
Barium	200	2.0	92.0	A								
Beryllium	5.0	1.0	2.8	L								
Cadmium	5.0	3.0	3.0	U								
Calcium	5000	17.0	159000	A								
Chromium	10.0	8.0	84.8	J								
Cobalt	50.0	9.0	33.8	A								
Copper	25.0	4.0	12.4	J								
Iron	100	25.0	44300	J								
Lead	3.0	2.0	4.0	J								
Magnesium	5000	48.0	102000	J								
Manganese	15.0	1.0	1050	J								
Mercury	0.20	0.20	0.20	U								
Nickel	40.0	18.0	39.6	L								
Potassium	5000	1270	7080	A								
Selenium	50.0	30.0	30.0	R								
Silver	10.0	4.0	4.0	U								
Sodium	5000	123	7460	J								
Thallium	10.0	3.0	3.0	UL								
Vanadium	50.0	6.0	99.0	A								
Zinc	20.0	4.0	278	J								
Cyanide			Not Analyzed									

Notes: All units are ug/l
Sample Quantitation Limit (SOL) is the Instrument Detection Limit (IDL).
The CRDLs and IDLs have been adjusted for dilution factors and percent solids as appropriate.

Qualifiers.....
 A - Acceptable (Quantitative) Data
 J - Estimated (Semi-quantitative) Data
 K - Biased High (Semi-quantitative) Data
 L - Biased Low (Semi-quantitative) Data
 B - Present in associated blanks (Semi-quantitative) Data
 UJ - Not Detected/Estimated Data
 U - Not Detected/Quantitative Data
 UL - Not Detected/Biased Low
 N - Tentative Identification
 R - Unusable Data
 Q - Not Applicable

AR302832

DISSOLVED (FILTERED) METALS

Phase 1A

Phase 1B

AR302833

October 1993

913-6773

SUMMARY OF CLP DISSOLVED INORGANICS ANALYSES - PHASE 1A
 BERKS LANDFILL REMEDIAL INVESTIGATION - PHASE 1A
 BERKS COUNTY, PENNSYLVANIA

Matrix: Groundwater Well

Parameter	Sample Point GR180/1A			Sample Point GVG01/1A			Sample Point GVG02/1A			Sample Point GVG03/1A		
	CRDL	IDL	Result	Qual	CRDL	IDL	Result	Qual	CRDL	IDL	Result	Qual
Aluminum	200	14.0	14.0	U	200	14.0	22.5	B	200	14.0	28.9	B
Antimony	60.0	14.0	14.0	U	60.0	14.0	14.0	U	60.0	14.0	14.0	U
Arsenic	10.0	3.0	4.8	A	10.0	3.0	3.0	U	10.0	3.0	3.0	U
Barium	200	0.40	30.7	B	200	0.40	39.6	A	200	0.40	11.6	A
Beryllium	5.0	0.30	0.30	U	5.0	0.30	0.30	U	5.0	0.30	0.30	U
Cadmium	5.0	1.9	2.3	B	5.0	1.9	1.9	UL	5.0	1.9	1.9	UL
Calcium	5000	4.1	52200	A	5000	4.1	110000	J	5000	4.1	29400	J
Chromium	10.0	3.6	3.6	U	10.0	3.6	3.6	UL	10.0	3.6	3.6	UL
Cobalt	50.0	2.0	2.0	B	50.0	2.0	2.0	U	50.0	2.0	2.5	B
Copper	25.0	1.9	6.7	B	25.0	1.9	3.9	B	25.0	1.9	3.5	B
Iron	100	9.2	9.2	UL	100	9.2	56.7	B	100	9.2	49.1	B
Lead	3.0	1.0	1.1	B	3.0	1.0	1.0	U	3.0	1.0	1.0	B
Magnesium	5000	12.2	8930	A	5000	12.2	27800	J	5000	12.2	9340	J
Manganese	15.0	0.60	15.2	A	15.0	0.60	857	J	15.0	0.60	51.7	J
Mercury	0.20	0.05	0.05	A	0.20	0.04	0.21	B	0.20	0.04	0.21	B
Nickel	40.0	7.9	7.9	U	40.0	7.9	12.3	B	40.0	7.9	12.0	B
Potassium	5000	64.0	6030	A	5000	64.0	2710	A	5000	64.0	976	A
Selenium	5.0	5.0	5.0	U	5.0	5.0	5.0	U	5.0	5.0	5.0	U
Silver	10.0	2.0	2.0	U	10.0	2.0	2.0	UL	10.0	2.0	2.0	UL
Sodium	5000	8.0	15900	A	5000	8.0	9850	J	5000	8.0	9300	J
Thallium	10.0	2.0	2.0	U	10.0	2.0	2.0	U	10.0	2.0	2.0	U
Venadium	50.0	1.8	1.8	U	50.0	1.8	1.8	U	50.0	1.8	5.0	U
Zinc	20.0	1.2	7.4	B	20.0	1.2	26.3	B	20.0	1.2	5.5	B

Notes: All units are ug/l

Sample Quantitation Limit (SQL) is the Instrument Detection Limit (IDL).

The CRDLs and IDLs have been adjusted for dilution factors and percent solids as appropriate.

Qualifiers.....
 A - Acceptable (Quantitative) Data
 J - Estimated (Semi-quantitative) Data
 K - Biased High (Semi-quantitative) Data
 L - Biased Low (Semi-quantitative) Data
 B - Present in associated blanks (Semi-quantitative) Data
 UJ - Not Detected/Estimated Data
 U - Not Detected/Quantitative Data
 UL - Not Detected/Biased Low
 N - Tentative Identification
 R - Unusable Data
 Q - Not Applicable

AR302834

SUMMARY OF CLP DISSOLVED INORGANICS ANALYSES - PHASE 1A

BERKS LANDFILL REMEDIAL INVESTIGATION - PHASE 1A
BERKS COUNTY, PENNSYLVANIA

Matrix: Groundwater Well

Parameter	Sample Point GWC3D/1A			Sample Point GWC3S/1A			Sample Point GJ-C5/1A			Sample Point GJ-16/1A		
	CRDL	IDL	Result	Qual	CRDL	IDL	Result	Qual	CRDL	IDL	Result	Qual
Aluminum	200	14.0	14.0	U	200	14.0	17.0	B	200	14.0	14.0	B
Antimony	60.0	14.0	14.0	U	60.0	14.0	14.0	U	60.0	14.0	14.0	U
Arsenic	10.0	3.0	3.0	U	10.0	3.0	3.9	A	10.0	3.0	3.0	U
Barium	200	0.40	50.7	A	200	0.40	80.4	A	200	0.40	32.2	A
Beryllium	5.0	0.30	0.30	U	5.0	0.30	0.30	U	5.0	0.30	0.30	U
Cadmium	5.0	1.9	1.9	U	5.0	1.9	1.9	U	5.0	1.9	1.9	U
Calcium	5000	4.1	168000	J	5000	4.1	216000	J	5000	4.1	91300	J
Chromium	10.0	3.6	3.6	UL	10.0	3.6	3.6	UL	10.0	3.6	3.6	UL
Cobalt	50.0	2.0	2.0	U	50.0	2.0	13.2	A	50.0	2.0	2.0	U
Copper	25.0	1.9	1.9	U	25.0	1.9	9.9	B	25.0	1.9	1.9	U
Iron	100	9.2	12000	A	100	9.2	580	L	100	9.2	4490	A
Lead	3.0	1.0	1.0	U	3.0	1.0	3.0	B	3.0	1.0	1.0	U
Magnesium	5000	12.2	79600	J	5000	12.2	133000	J	5000	12.2	36700	J
Manganese	15.0	0.60	408	J	15.0	0.60	8460	J	15.0	0.60	559	J
Mercury	0.20	0.04	0.04	UL	0.20	0.04	0.04	UL	0.20	0.04	0.04	UL
Nickel	40.0	7.9	7.9	U	40.0	7.9	90.5	A	40.0	7.9	10.9	B
Potassium	5000	64.0	1810	A	5000	64.0	3360	A	5000	64.0	5280	A
Selenium	5.0	5.0	2.2	B	5.0	5.0	5.0	U	5.0	5.0	5.0	U
Silver	10.0	2.0	2.0	B	10.0	2.0	2.0	U	10.0	2.0	3.0	B
Sodium	5000	8.0	13500	J	5000	8.0	42400	J	5000	8.0	22800	J
Thallium	10.0	2.0	2.0	UL	10.0	2.0	2.0	UL	10.0	2.0	2.0	UL
Vanadium	50.0	1.8	1.8	U	50.0	1.8	1.8	U	50.0	1.8	1.8	U
Zinc	20.0	1.2	16.4	B	20.0	1.2	44.9	A	20.0	1.2	8.8	B

Notes: All units are ug/l
Sample Quantitation Limit (SQL) is the Instrument Detection Limit (IDL).
The CRDLs and IDLs have been adjusted for dilution factors and percent solids as appropriate.

Qualifiers.....
A - Acceptable (Quantitative) Data
J - Estimated (Semi-quantitative) Data
K - Biased High (Semi-quantitative) Data
L - Biased Low (Semi-quantitative) Data
B - Present in associated blanks (Semi-quantitative) Data
UJ - Not Detected/Estimated Data
UL - Not Detected/Quantitative Data
UL - Not Detected/Biased Low
N - Tentative Identification
R - Unusable Data
Q - Not Applicable

AR302835

October 1993

SUMMARY OF CLP DISSOLVED INORGANICS ANALYSES - PHASE 1A
 BERKS LANDFILL REMEDIAL INVESTIGATION - PHASE 1A
 BERKS COUNTY, PENNSYLVANIA

913-6773

Matrix: Groundwater Well

Parameter	Sample Point GV18S/1A Date Sampled: 09/03/92 Dilution Factor: 1.0 Percent Solids: NA %			Sample Point GV18S/FO Date Sampled: 09/03/92 Dilution Factor: 1.0 Percent Solids: NA %			Sample Point Date Sampled: Dilution Factor: Percent Solids: %			Sample Point Date Sampled: Dilution Factor: Percent Solids: %		
	CRDL	IDL	Result	Qual	CRDL	IDL	Result	Qual	CRDL	IDL	Result	Qual
Aluminum	200	14.0	59.1	B	200	14.0	25.5	B				
Antimony	60.0	14.0	14.0	U	60.0	14.0	14.0	U				
Arsenic	10.0	3.0	3.0	U	10.0	3.0	3.0	U				
Barium	200	0.40	61.0	A	200	0.40	63.5	A				
Beryllium	5.0	0.30	0.30	U	5.0	0.30	0.30	U				
Cadmium	5.0	1.9	1.9	U	5.0	1.9	1.9	U				
Calcium	5000	4.1	216000	J	5000	4.1	221000	J				
Chromium	10.0	3.6	3.6	UL	10.0	3.6	3.6	UL				
Cobalt	50.0	2.0	2.0	U	50.0	2.0	2.0	U				
Copper	25.0	1.9	1.9	U	25.0	1.9	1.9	U				
Iron	100	9.2	10.4	B	100	9.2	9.2	UL				
Lead	3.0	1.0	1.2	B	3.0	1.0	1.0	U				
Magnesium	5000	12.2	50900	J	5000	12.2	52500	J				
Manganese	15.0	0.60	72.1	J	15.0	0.60	69.8	J				
Mercury	0.20	0.04	0.04	UL	0.20	0.04	0.04	UL				
Nickel	40.0	7.9	7.9	U	40.0	7.9	7.9	U				
Potassium	5000	64.0	2630	A	5000	64.0	2790	A				
Selenium	5.0	5.0	5.0	U	5.0	5.0	5.0	U				
Silver	10.0	2.0	3.6	B	10.0	2.0	3.6	B				
Sodium	5000	8.0	16100	J	5000	8.0	16700	J				
Thallium	10.0	2.0	2.0	UL	10.0	2.0	2.0	UL				
Vanadium	50.0	1.8	1.8	U	50.0	1.8	1.8	U				
Zinc	20.0	1.2	11.3	B	20.0	1.2	15.2	B				

Notes: All units are ug/l
 Sample Quantitation Limit (SQL) is the Instrument Detection Limit (IDL).
 The CRDLs and IDLs have been adjusted for dilution factors and percent solids as appropriate.

Qualifiers.....
 A - Acceptable (Quantitative) Data
 J - Estimated (Semi-quantitative) Data
 K - Biased High (Semi-quantitative) Data
 L - Biased Low (Semi-quantitative) Data
 B - Present in associated blanks (Semi-quantitative) Data
 UJ - Not Detected/Estimated Data
 U - Not Detected/Quantitative Data
 UL - Not Detected/Biased Low
 M - Tentative Identification
 R - Unusable Data
 Q - Not Applicable

AR302836

October 1993

913-6773

SUMMARY OF CLP DISSOLVED INORGANICS ANALYSES - PHASE 1B

BERKS LANDFILL REMEDIAL INVESTIGATION - PHASE 1B
BERKS COUNTY, PENNSYLVANIA

Matrix: Groundwater

Parameter	Sample Point GUC01/1B			Sample Point GUC02/1B			Sample Point GUC03/1B			Sample Point GUC03/1B		
	CRDL	IDL	Result	Qual	CRDL	IDL	Result	Qual	CRDL	IDL	Result	Qual
Aluminum	200	44	57.4	B	200	44	59.5	B	200	35	35.0	U
Antimony	60	35	35.0	U	60	35	35.0	U	60	36	36.0	U
Arsenic	10	4	4.0	U	10	4	4.0	U	10	4	4.0	UJ
Barium	200	2	2.0	U	200	2	15.7	A	200	2	18.9	B
Beryllium	5	1	1.0	U	5	1	1.0	U	5	1	1.0	U
Cadmium	5	3	3.0	U	5	3	3.0	U	5	5	5.0	U
Calcium	5000	17	29500	A	5000	17	33100	A	5000	24	146000	A
Chromium	10	8	8.0	U	10	8	8.0	U	10	8	8.0	U
Cobalt	50	9	9.0	U	50	9	9.0	U	50	8	8.0	U
Copper	25	4	4.0	U	25	4	4.0	U	25	4	4.0	U
Iron	100	25	66.8	A	100	25	53.5	A	100	41	41.0	U
Lead	3	2	2.0	UL	3	2	2.0	UL	3	2	2.0	UL
Magnesium	5000	48	10800	A	5000	48	11800	A	5000	59	49700	A
Manganese	15	1	502	A	15	1	28.9	B	15	1	30.6	B
Mercury	0.2	0.2	0.20	UL	0.2	0.2	0.20	UL	0.2	0.2	0.20	UL
Nickel	40	18	18.0	U	40	18	18.0	U	40	17	17.0	U
Potassium	5000	1270	1270	U	5000	1270	1270	U	5000	1352	1370	A
Selenium	5	3	3.0	UL	5	3	3.0	UL	50	30	30.0	UJ
Silver	10	4	4.0	UL	10	4	4.0	UL	10	7	7.0	UL
Sodium	5000	123	7520	A	5000	123	9950	A	5000	251	12600	A
Thallium	10	3	3.0	R	10	3	3.0	R	100	40	40.0	R
Vanadium	50	6	6.0	U	50	6	6.0	U	50	6	6.0	U
Zinc	20	4	4.0	U	20	4	6.0	B	20	4	16.7	B
Cyanide	10	10	Not Analyzed		10	10	Not Analyzed		10	10	Not Analyzed	

Notes: All units are ug/l
 Sample Quantitation Limit (SQL) is the Instrument Detection Limit (IDL).
 The CRDLs and IDLs have been adjusted for dilution factors and percent solids as appropriate.

Qualifiers.....
 A - Acceptable (Quantitative) Data
 J - Estimated (Semi-quantitative) Data
 K - Biased High (Semi-quantitative) Data
 L - Biased Low (Semi-quantitative) Data
 B - Present in associated blanks (Semi-quantitative) Data
 UJ - Not Detected/Estimated Data
 U - Not Detected/Quantitative Data
 UL - Not Detected/Biased Low
 N - Tentative Identification
 R - Unusable Data
 O - Not Applicable

AR302837

SUMMARY OF CLP DISSOLVED INORGANICS ANALYSES - PHASE 1B
 BERKS LANDFILL REMEDIAL INVESTIGATION - PHASE 1B
 BERKS COUNTY, PENNSYLVANIA

Matrix: Groundwater

Parameter	Sample Point GMC03S/1B			Sample Point GMC040/1B			Sample Point GMC04S/1B			Sample Point GMC060/1B		
	CRDL	IDL	Qual	CRDL	IDL	Qual	CRDL	IDL	Qual	CRDL	IDL	Qual
Aluminum	200	35	U	200	35	B	200	35	U	200	44	U
Antimony	60	36	U	60	36	U	60	36	U	60	35	U
Arsenic	10	4	J	10	3	A	10	3	A	10	4	U
Barium	200	2	A	200	2	A	200	2	A	200	2	A
Beryllium	5	1	A	5	1	U	5	1	U	5	1	U
Cadmium	5	5	U	5	5	U	5	5	U	5	3	U
Calcium	5000	24	A	5000	24	U	5000	24	U	5000	17	U
Chromium	10	8	A	10	8	U	10	8	U	10	8	U
Chromium	50	8	A	50	8	U	50	8	U	50	9	U
Cobalt	25	4	A	25	4	U	25	4	U	25	4	U
Copper	100	41	U	100	41	A	100	41	U	100	25	U
Iron	3	2	UL	3	2	UL	3	2	UL	3	2	UL
Lead	5000	59	A	5000	59	U	5000	59	A	5000	48	A
Manganese	15	1	A	15	1	A	15	1	A	15	1	A
Mercury	0.2	0.2	UL	0.2	0.2	UL	0.2	0.2	UL	0.2	0.2	UL
Nickel	40	17	A	40	17	U	40	17	U	40	18	A
Potassium	5000	1352	A	5000	1352	U	5000	1352	A	5000	1270	A
Selenium	5	3	UJ	5	30	R	5	30	U	5	3	UL
Silver	10	7	UL	10	7	UL	10	7	UL	10	4	UL
Sodium	5000	251	A	5000	251	U	5000	251	A	5000	123	A
Thallium	100	40	R	100	40	R	100	40	R	100	3	R
Vanadium	50	6	U	50	6	U	50	6	U	50	6	R
Zinc	20	4	U	20	4	U	20	4	U	20	4	A
Cyanide	10	10	B	10	10	B	10	10	B	10	10	B
			Not Analyzed			Not Analyzed			Not Analyzed			Not Analyzed
			35.0			396			35.0			104
			36.0			36.0			36.0			35.0
			10.2			7.3			10.2			4.0
			87.3			173			398			60.3
			1.0			1.0			1.0			1.0
			5.0			5.0			5.0			3.0
			230000			16600			148000			33900
			8.0			8.0			8.0			8.0
			12.4			8.0			8.0			9.0
			4.0			4.0			4.0			4.0
			1030			182			959			25.0
			2.0			2.0			2.0			2.0
			147000			3140			47000			11300
			8320			78.3			970			4.1
			0.20			0.20			0.20			0.20
			75.3			17.0			4370			18.0
			2640			1670			30.0			3.0
			3.0			30.0			7.0			4.0
			39000			21000			17900			8180
			40.0			4.0			40.0			3.0
			6.0			44.7			6.0			9.5
			15.2			Not Analyzed			80.0			4.4

Notes: All units are ug/L
 Sample Quantitation Limit (SQL) is the Instrument Detection Limit (IDL).
 The CRDLs and IDLs have been adjusted for dilution factors and percent solids as appropriate.

Qualifiers.....
 A - Acceptable (Quantitative) Data
 J - Estimated (Semiquantitative) Data
 K - Biased High (Semiquantitative) Data
 L - Biased Low (Semiquantitative) Data
 B - Present in associated blanks (Semiquantitative) Data
 UJ - Not Detected/Estimated Data
 U - Not Detected/Quantitative Data
 UL - Not Detected/Biased Low
 M - Tentative Identification
 R - Unusable Data
 Q - Not Applicable

SUMMARY OF CLP DISSOLVED INORGANICS ANALYSES - PHASE 1B

BERKS LANDFILL REMEDIAL INVESTIGATION - PHASE 1B
BERKS COUNTY, PENNSYLVANIA

Matrix: Groundwater

Parameter	Sample Point GUC06S/1B			Sample Point GUC07D/1B			Sample Point GUC07S/1B			Sample Point GUC02/1B		
	CRDL	IDL	Result	Qual	CRDL	IDL	Result	Qual	CRDL	IDL	Result	Qual
Aluminum	200	44	58.9	B	200	44	54.8	B	200	44	61.5	B
Antimony	60	35	35.0	U	60	35	35.0	U	60	35	35.0	U
Arsenic	10	4	4.0	U	10	4	4.0	U	10	4	4.0	U
Barium	200	2	97.7	A	200	2	45.0	A	200	2	334	A
Beryllium	5	1	1.0	U	5	1	1.0	U	5	1	1.0	U
Cadmium	5	3	3.0	U	5	3	3.0	U	5	3	3.0	U
Calcium	5000	17	21800	A	5000	17	33600	A	5000	17	325000	A
Chromium	10	8	8.0	U	10	8	8.0	U	10	8	8.0	U
Chromium	50	9	9.0	U	50	9	9.0	U	50	9	9.0	U
Cobalt	25	4	4.0	U	25	4	4.0	U	25	4	4.0	U
Copper	100	25	26.8	U	100	25	25.0	U	100	25	26.5	A
Iron	3	2	2.0	U	3	2	2.0	U	3	2	2.0	U
Lead	5000	48	7050	UL	5000	48	7100	A	5000	48	68500	A
Magnesium	15	1	11.0	B	15	1	30.1	B	15	1	1330	A
Manganese	0.2	0.2	0.20	UL	0.2	0.2	0.20	UL	0.2	0.2	0.20	UL
Mercury	40	18	18.0	U	40	18	18.0	U	40	18	29.4	A
Nickel	5000	1270	1610	A	5000	1270	1660	A	5000	1270	4300	A
Potassium	5	3	3.0	UL	5	3	3.0	UL	5	3	3.0	UL
Selenium	10	4	4.0	U	10	4	4.0	U	10	4	4.0	U
Silver	5000	123	7290	A	5000	123	10500	A	5000	123	24800	A
Sodium	10	3	3.0	R	10	3	3.0	R	100	30	30.0	R
Thallium	50	6	6.0	U	50	6	6.0	U	50	6	6.0	U
Vanadium	20	4	4.0	U	20	4	4.0	U	20	4	22.1	B
Zinc	10	10	Not Analyzed	U	10	10	Not Analyzed	U	10	10	Not Analyzed	B
Cyanide												

Notes: All units are ug/l
Sample Quantitation Limit (SQL) is the Instrument Detection Limit (IDL).
The CRDLs and IDLs have been adjusted for dilution factors and percent solids as appropriate.

Qualifiers.....
A - Acceptable (Quantitative) Data
J - Estimated (Semi-quantitative) Data
K - Biased High (Semi-quantitative) Data
L - Biased Low (Semi-quantitative) Data
B - Present in associated blanks (Semi-quantitative) Data
UL - Not Detected/Estimated Data
UJ - Not Detected/Quantitative Data
U - Not Detected/Quantitative Data
UL - Not Detected/Biased Low
N - Tentative Identification
R - Unusable Data
Q - Not Applicable

October 1993

SUMMARY OF CLP DISSOLVED INORGANICS ANALYSES - PHASE 1B

913-6773

BERKS LANDFILL REMEDIAL INVESTIGATION - PHASE 1B
BERKS COUNTY, PENNSYLVANIA

Matrix: Groundwater

Parameter	Sample Point GMG01/1B Date Sampled: 06/30/93 Dilution Factor: 1.0 % Percent Solids: NA			Sample Point GMG02/1B Date Sampled: 06/30/93 Dilution Factor: 1.0 % Percent Solids: NA			Sample Point GMG03/1B Date Sampled: 06/30/93 Dilution Factor: 1.0 % Percent Solids: NA			Sample Point GMG04/1B Date Sampled: 07/09/93 Dilution Factor: 1.0 % Percent Solids: NA		
	CRDL	IDL	Result	Qual	CRDL	IDL	Result	Qual	CRDL	IDL	Result	Qual
Aluminum	200	35	35.0	U	200	35	35.0	U	200	44	56.1	B
Antimony	60	36	36.0	U	60	36	36.0	U	60	35	35.0	U
Arsenic	10	4	4.0	UJ	10	4	4.0	UJ	10	4	18.1	A
Barium	200	2	26.1	B	200	2	11.1	B	200	2	76.0	U
Beryllium	5	1	1.0	U	5	1	1.0	U	5	1	1.0	U
Cadmium	5	5	5.0	U	5	5	5.0	U	5	3	3.0	U
Calcium	5000	24	23600	A	5000	24	31600	A	5000	17	247000	U
Chromium	10	8	8.0	U	10	8	8.0	U	10	8	8.0	U
Cobalt	50	8	8.0	U	50	8	8.0	U	50	9	9.0	U
Copper	25	4	4.0	U	25	4	4.0	U	25	4	4.0	U
Iron	100	41	41.0	U	100	41	41.0	U	100	25	762	A
Lead	3	2	2.0	UL	3	2	2.0	UL	3	1	2.6	L
Magnesium	5000	59	13200	A	5000	59	9570	A	5000	48	42900	A
Manganese	15	1	33.1	A	15	1	6.8	B	15	1	1590	A
Mercury	0.2	0.2	0.20	UL	0.2	0.2	0.20	UL	0.2	0.2	0.20	UL
Nickel	40	17	17.0	U	40	17	17.0	U	40	18	19.7	A
Potassium	5000	1352	1350	U	5000	1352	1350	U	5000	1270	3640	A
Selenium	5	3	3.0	UJ	5	3	30.0	UJ	5	3	3.0	UL
Silver	10	7	7.0	UL	10	7	7.0	UL	10	4	4.0	UL
Sodium	5000	251	13100	A	5000	251	9920	A	5000	123	40400	A
Thallium	10	4	4.0	R	10	4	4.0	R	10	30	30.0	R
Vanadium	50	6	6.0	U	50	6	6.0	U	50	6	6.0	U
Zinc	20	4	7.9	B	20	4	8.2	B	20	4	37.0	U
Cyanide	10	10	Not Analyzed	B	10	10	Not Analyzed	B	10	10	Not Analyzed	B

Notes: All units are ug/l
Sample Quantitation Limit (SQL) is the Instrument Detection Limit (IDL).
The CRDLs and IDLs have been adjusted for dilution factors and percent solids as appropriate.

Qualifiers.....
A - Acceptable (Quantitative) Data
J - Estimated (Semi-quantitative) Data
K - Biased High (Semi-quantitative) Data
L - Biased Low (Semi-quantitative) Data
B - Present in associated blanks (Semi-quantitative) Data
UJ - Not Detected/Estimated Data
U - Not Detected/Quantitative Data
UL - Not Detected/Biased Low
M - Tentative Identification
R - Unusable Data
Q - Not Applicable

AR302840

SUMMARY OF CLP DISSOLVED INORGANICS ANALYSES - PHASE 1B

BERKS LANDFILL REMEDIAL INVESTIGATION - PHASE 1B
BERKS COUNTY, PENNSYLVANIA

Matrix: Groundwater

Parameter	Sample Point GUG05/1B			Sample Point GUG06/1B			Sample Point GUG07/1B			Sample Point GUG08/1B						
	CRDL	IDL	Result	Qual	CRDL	IDL	Result	Qual	CRDL	IDL	Result	Qual	CRDL	IDL	Result	Qual
Aluminum	200	35	43.4	B	200	35	35.0	U	200	35	59.9	B	200	35	35.0	U
Antimony	60	36	36.0	U	60	36	36.0	U	60	36	36.0	U	60	36	36.0	U
Arsenic	10	4	11.6	A	10	4	4.0	U	10	4	7.3	A	10	4	4.0	U
Barium	200	2	148	A	200	2	2.0	U	200	2	7.5	B	200	2	15.7	B
Beryllium	5	1	1.0	U	5	1	1.0	U	5	1	1.0	U	5	1	1.0	U
Cadmium	5	5	5.0	U	5	5	5.0	U	5	5	5.0	U	5	5	5.0	U
Calcium	5000	24	232000	A	5000	24	40400	A	5000	24	22600	A	5000	24	28900	A
Chromium	10	8	8.0	U	10	8	8.0	U	10	8	8.0	U	10	8	8.0	U
Cobalt	50	8	8.0	U	50	8	8.0	U	50	8	8.0	U	50	8	8.0	U
Copper	25	4	4.0	U	25	4	4.0	U	25	4	4.0	U	25	4	4.0	U
Iron	100	41	128	A	100	41	41.0	U	100	41	162	A	100	41	41.0	U
Lead	3	2	2.0	UL	3	2	2.0	UL	3	2	2.0	UL	3	2	2.0	UL
Magnesium	5000	59	48900	A	5000	59	14900	A	5000	59	1660	A	5000	59	5640	A
Manganese	15	1	1070	A	15	1	2.5	B	15	1	8.5	B	15	1	361	A
Mercury	0.2	0.2	0.20	UL	0.2	0.2	0.20	UL	0.2	0.2	0.20	UL	0.2	0.2	0.20	UL
Nickel	40	17	20.4	A	40	17	17.0	U	40	17	17.0	U	40	17	17.0	U
Potassium	5000	1352	5090	A	5000	1352	1350	U	5000	1352	1350	U	5000	1352	1350	U
Selenium	50	30	30.0	UJ	50	30	30.0	UJ	50	30	3.0	UJ	50	30	3.0	UJ
Silver	10	7	7.0	UL	10	7	7.0	UL	10	7	7.0	UL	10	7	7.0	UL
Sodium	5000	251	66600	A	5000	251	7240	A	5000	251	11700	A	5000	251	5570	A
Thallium	100	40	40.0	R	100	40	4.0	R	100	40	4.0	R	100	40	4.0	R
Vanadium	50	6	6.0	U	50	6	6.3	A	50	6	12.0	A	50	6	6.0	U
Zinc	20	4	20.4	B	20	4	9.7	B	20	4	4.0	U	20	4	6.7	B
Cyanide	10	10	Not Analyzed		10	10	Not Analyzed		10	10	Not Analyzed		10	10	Not Analyzed	

Notes: All units are ug/l
Sample Quantitation Limit (SQL) is the Instrument Detection Limit (IDL).
The CRDLs and IDLs have been adjusted for dilution factors and percent solids as appropriate.

Qualifiers.....
A - Acceptable (Quantitative) Data
J - Estimated (Semi-quantitative) Data
K - Biased High (Semi-quantitative) Data
L - Biased Low (Semi-quantitative) Data
B - Present in associated blanks (Semi-quantitative) Data
.....
UJ - Not Detected/Estimated Data
UL - Not Detected/Quantitative Data
UL - Not Detected/Biased Low
.....
M - Tentative Identification
R - Unusable Data
Q - Not Applicable

AR302841

October 1993

SUMMARY OF CLP DISSOLVED INORGANICS ANALYSES - PHASE 1B

913-6773

BERKS LANDFILL REMEDIAL INVESTIGATION - PHASE 1B
BERKS COUNTY, PENNSYLVANIA

Matrix: Groundwater

Parameter	Sample Point GUG10/1B				Sample Point GUG11/1B				Sample Point GUG11/1B				Sample Point GUG12/1B			
	CRDL	IDL	Result	Qual	CRDL	IDL	Result	Qual	CRDL	IDL	Result	Qual	CRDL	IDL	Result	Qual
Aluminum	200	35	35.0	U	200	44	51.7	B	200	44	44.0	U	200	35	35.0	U
Antimony	60	36	36.0	U	60	35	35.0	U	60	35	42.1	A	60	36	36.0	U
Arsenic	10	4	4.0	U	10	4	4.0	U	10	4	4.0	U	10	4	4.2	A
Barium	200	2	6.3	B	200	2	1430	A	200	2	1420	A	200	2	475	A
Beryllium	5	1	1.0	U	5	1	1.0	U	5	1	1.0	U	5	1	1.0	U
Cadmium	5	5	5.0	U	5	3	3.0	U	5	3	3.0	U	5	5	5.0	U
Calcium	5000	24	23000	A	5000	17	470000	A	5000	17	470000	A	5000	24	430000	A
Chromium	10	8	8.0	U	10	8	8.0	U	10	8	8.0	U	10	8	8.0	U
Cobalt	50	8	8.0	U	50	9	9.0	U	50	9	9.0	U	50	8	8.0	U
Copper	25	4	4.0	U	25	4	4.0	U	25	4	4.0	U	25	4	4.0	U
Iron	100	41	41.0	U	100	25	481	A	100	25	481	A	100	25	182	A
Lead	3	2	2.0	UL	3	1	2.0	UL	3	1	2.0	UL	3	2	11.2	L
Magnesium	5000	59	7180	A	5000	48	108000	A	5000	48	108000	A	5000	59	99100	A
Manganese	15	1	36.9	A	15	1	5110	A	15	1	5110	A	15	1	2470	A
Mercury	0.2	0.2	0.20	UL	0.2	0.2	0.20	UL	0.2	0.2	0.20	UL	0.2	0.2	0.20	UL
Nickel	40	17	17.0	U	40	18	47.5	A	40	18	47.4	A	40	17	24.1	A
Potassium	5000	1352	1350	U	5000	1270	8840	A	5000	1270	9080	A	5000	1352	7260	A
Selenium	5	3	3.0	UJ	5	3	3.0	UL	5	3	3.0	UL	5	30	30.0	UJ
Silver	10	7	7.0	UL	10	4	4.0	U	10	4	4.0	U	10	7	7.0	UL
Sodium	5000	251	7580	R	5000	123	155000	A	5000	123	155000	A	5000	251	27800	R
Thallium	10	4	4.0	R	10	30	30.0	R	10	30	30.0	R	10	40	40.0	R
Vanadium	50	6	6.0	U	50	6	6.0	U	50	6	6.0	U	50	6	6.0	U
Zinc	20	4	4.0	U	20	4	12.6	B	20	4	11.1	B	20	4	4.0	U
Cyanide	10	10	Not Analyzed	U	10	10	Not Analyzed	B	10	10	Not Analyzed	B	10	10	Not Analyzed	U

Notes: All units are ug/l

Sample Quantitation Limit (SQL) is the Instrument Detection Limit (IDL)

The CRDLs and IDLs have been adjusted for dilution factors and percent solids as appropriate.

Qualifiers:
 A - Acceptable (Quantitative) Data
 J - Estimated (Semi-quantitative) Data
 K - Biased High (Semi-quantitative) Data
 L - Biased Low (Semi-quantitative) Data
 B - Present in associated blanks (Semi-quantitative) Data
 UJ - Not Detected/Estimated Data
 U - Not Detected/Quantitative Data
 UL - Not Detected/Biased Low
 N - Tentative Identification
 R - Unusable Data
 Q - Not Applicable

AR302842

SUMMARY OF CLP DISSOLVED INORGANICS ANALYSES - PHASE 1B
 BERKS LANDFILL REMEDIAL INVESTIGATION - PHASE 1B
 BERKS COUNTY, PENNSYLVANIA

Matrix: Groundwater

Parameter	Sample Point G4G12/1D			Sample Point G4G13/1B			Sample Point G4H03/1B			Sample Point G4H11/1B		
	CRDL	IDL	Result	Qual	CRDL	IDL	Result	Qual	CRDL	IDL	Result	Qual
Aluminum	200	35	35.0	U	200	44	49.4	B	200	44	44.0	U
Antimony	60	36	36.0	U	60	35	38.2	A	60	35	35.0	U
Arsenic	10	4	4.6	J	10	4	14.6	A	10	4	9.7	A
Barium	200	2	480	A	200	2	11.1	A	200	2	123	A
Beryllium	5	1	1.0	U	5	1	1.0	U	5	1	1.0	U
Cadmium	5	5	5.0	U	5	3	3.0	U	5	5	5.0	U
Calcium	5000	24	432000	A	5000	17	46200	A	5000	17	160000	A
Chromium	10	8	8.0	U	10	8	8.0	U	10	8	8.0	U
Chromium	50	8	8.0	U	50	9	9.0	U	50	8	8.0	U
Cobalt	25	4	4.0	U	25	4	4.0	U	25	4	4.0	U
Copper	100	41	168	A	100	25	25.0	U	100	25	3640	A
Iron	3	2	14.0	L	3	1	2.0	U	3	1	2.2	L
Lead	5000	59	100000	A	5000	48	13400	A	5000	48	103000	A
Magnesium	15	1	2480	A	15	1	21.2	B	15	1	13900	A
Manganese	0.2	0.2	0.20	UL	0.2	0.2	0.20	UL	0.2	0.2	0.20	UL
Mercury	40	17	21.2	UL	40	18	18.0	UL	40	18	77.0	UL
Nickel	5000	1552	7340	A	5000	1270	1270	U	5000	1270	3950	A
Potassium	50	30	30.0	UJ	5	3	3.0	UL	5	3	3.0	UL
Selenium	10	7	7.0	UL	10	4	4.0	UL	10	4	4.0	UL
Silver	5000	251	28400	A	5000	123	12500	UL	5000	123	196000	A
Sodium	100	40	40.0	R	10	3	3.0	R	100	30	30.0	R
Thallium	50	6	6.0	U	50	6	12.3	A	50	6	6.0	U
Vanadium	20	4	4.0	U	20	4	4.0	U	20	4	20.6	U
Zinc	10	10	Not Analyzed	U	10	10	Not Analyzed	U	10	10	Not Analyzed	B
Cyanide												

Notes: All units are ug/l.
 Sample quantitation limit (SQL) is the Instrument Detection Limit (IDL).
 The CRDLs and IDLs have been adjusted for dilution factors and percent solids as appropriate.

Qualifiers.....
 A - Acceptable (Quantitative) Data
 J - Estimated (Semi-quantitative) Data
 K - Biased High (Semi-quantitative) Data
 L - Biased Low (Semi-quantitative) Data
 B - Present in associated blanks (Semi-quantitative) Data
 UJ - Not Detected/Estimated Data
 U - Not Detected/Quantitative Data
 UL - Not Detected/Biased Low
 M - Tentative Identification
 R - Unusable Data
 O - Not Applicable

AR302843

SUMMARY OF CLP DISSOLVED INORGANICS ANALYSES - PHASE 1B
 BERKS LANDFILL REMEDIAL INVESTIGATION - PHASE 1B
 BERKS COUNTY, PENNSYLVANIA

Matrix: Groundwater

Parameter	Sample Point GUM16/1B			Sample Point GUM17/1B			Sample Point GUM19S/1B			Sample Point GUM190/1B		
	CRDL	IDL	Result	Qual	CRDL	IDL	Result	Qual	CRDL	IDL	Result	Qual
Aluminum	200	44	44.0	U	200	44	56.2	B	200	35	43.0	B
Antimony	60	35	38.0	A	60	35	35.0	U	60	36	36.0	U
Arsenic	10	4	9.0	A	10	4	4.0	U	10	3	3.0	U
Barium	200	2	359	A	200	2	3.2	A	200	2	67.8	A
Beryllium	5	1	1.0	U	5	1	1.0	U	5	1	1.0	U
Cadmium	5	3	3.0	U	5	3	3.0	U	5	5	5.0	U
Calcium	5000	17	400000	A	5000	17	13300	A	5000	24	427000	A
Chromium	10	8	8.0	U	10	8	8.0	U	10	8	8.0	U
Cobalt	50	9	10.3	A	50	9	9.0	U	50	8	8.0	U
Copper	25	4	4.0	U	25	4	4.0	U	25	4	4.0	U
Iron	100	25	1110	A	100	25	254	A	100	41	87.6	A
Lead	3	1	2.0	UL	3	1	2.0	UL	3	2	2.0	UL
Magnesium	5000	48	105000	A	5000	48	5560	A	5000	59	52300	A
Manganese	15	1	12300	A	15	1	251	A	15	1	724	A
Mercury	0.2	0.2	0.20	UL	0.2	0.2	0.20	UL	0.2	0.2	0.20	UL
Nickel	40	18	36.3	UL	40	18	18.0	UL	40	17	17.7	UL
Potassium	5000	1270	13400	A	5000	1270	1270	U	5000	1352	3090	A
Selenium	5	3	3.0	UL	5	3	3.0	UL	5	3	3.0	UL
Silver	10	4	4.0	UL	10	4	4.0	UL	10	7	7.0	UL
Sodium	5000	123	80500	A	5000	123	3700	B	5000	251	10600	A
Thallium	100	30	30.0	R	100	30	3.0	R	100	40	40.0	R
Vanadium	50	6	6.0	U	50	6	6.0	U	50	6	6.0	U
Zinc	20	4	11.7	B	20	4	5.1	B	20	4	14.0	B
Cyanide	10	10	Not Analyzed		10	10	Not Analyzed		10	10	Not Analyzed	

Notes: All units are ug/l
 Sample Quantitation Limit (SQL) is the Instrument Detection Limit (IDL).
 The CRDLs and IDLs have been adjusted for dilution factors and percent solids as appropriate.

Qualifiers:
 A - Acceptable (Quantitative) Data
 J - Estimated (Semi-quantitative) Data
 K - Biased High (Semi-quantitative) Data
 L - Biased Low (Semi-quantitative) Data
 B - Present in associated blanks (Semi-quantitative) Data

 UL - Not Detected/Estimated Data
 U - Not Detected/Quantitative Data
 UL - Not Detected/Biased Low

 N - Tentative Identification
 R - Unusable Data
 Q - Not Applicable

AR302845

WET CHEMISTRY

Phase 1A

AR302846

October 1993

913-6773

SUMMARY OF WET CHEMISTRY CONVENTIONAL ANALYSES - PHASE 1A

BERKS COUNTY, PENNSYLVANIA

Matrix: Groundwater

Parameter	Sample Point GR-160/1A			Sample Point GM-16/1A			Sample Point GM-16S/1A			Sample Point GM-16S/FD		
	RL	Result	Qual	RL	Result	Qual	RL	Result	Qual	RL	Result	Qual
Alkalinity - CaCO3	2.0	47.0	A	2.0	494	A	2.0	378	A	2.0	368	A
Ammonia - N	0.10	0.50	A	0.10	0.50	A	0.10	0.20	A	0.10	0.10	A
Biochemical Oxygen Dema	2.0	33.0	J	2.0	5.0	J	2.0	7.0	J	2.0	10.0	J
Chemical Oxygen Demand	5.0	87.0	A	5.0	104	A	5.0	19.0	B	5.0	21.0	B
Chloride	2.0	130	A	2.0	680	A	2.0	252	A	2.0	265	A
Fluoride	0.10	0.10	A	0.10	0.10	A	0.10	0.10	A	0.10	0.10	U
Hardness - CaCO3	2.0	183	A	2.0	1070	A	2.0	704	U	2.0	723	A
Hexavalent Chromium		Not Analyzed	A	0.10	0.10	U	0.10	0.10	U	0.10	0.10	U
Nitrate/Nitrite - N	0.02	0.05	B	0.02	0.61	J	0.02	0.07	B	0.02	0.16	B
Sulfate	5.0	19.0	J	5.0	38	J	5.0	53.0	J	5.0	30.0	B
Total Dissolved Solids	5.0	396	A	5.0	2050	A	5.0	1350	L	5.0	1430	L
Total Inorganic Carbon	1.0	9.0	A	1.0	130	A	1.0	94.0	A	1.0	100	A
Total Kjeldahl Nitrogen	0.10	0.69	B	0.10	2.6	B	0.10	0.40	B	0.10	0.40	B
Total Organic Carbon	1.0	11.8	A	1.0	30	A	1.0	4.4	A	1.0	6.3	A
Total Phosphate - P	0.05	0.20	B	0.05	0.21	B	0.05	0.05	U	0.05	0.05	B
Total Suspended Solids	5.0	25.0	A	5.0	163	A	5.0	21.0	U	5.0	16.0	L
Turbidity	0.20	22.0	J	0.20	118	J	0.20	4.8	L	0.20	2.3	J

Notes: All units are mg/l except Turbidity which is NTU.
 The Reporting Limits (RLs) have been adjusted for percent solids.
 The Qual column indicates the Qualifier applied to the result following data validation (see below).

- Qualifiers.....
- A - Acceptable (Quantitative) Data
 - J - Estimated (Semi-quantitative) Data
 - K - Biased High (Semi-quantitative) Data
 - L - Biased Low (Semi-quantitative) Data
 - B - Present in associated blanks (Semi-quantitative) Data
 - UJ - Not Detected/Estimated Data
 - U - Not Detected/Quantitative Data
 - UL - Not Detected/Biased Low
 - R - Unusable Data
 - Q - Not Applicable
 - N - Tentative Identification
 - N* - Tentative Identification (result below laboratory MDL)

AR302847

SUMMARY OF MET CHEMISTRY CONVENTIONAL ANALYSES - PHASE 1A

BERKS COUNTY, PENNSYLVANIA

Matrix: Groundwater

Parameter	Sample Point GWC-05/1A Date Sampled: 09/03/92 Dilution Factor: NA Percent Solids: NA %			Sample Point GWC-3D/1A Date Sampled: 09/04/92 Dilution Factor: NA Percent Solids: NA %			Sample Point GWC-3S/1A Date Sampled: 09/03/92 Dilution Factor: NA Percent Solids: NA %			Sample Point GWC-01/1A Date Sampled: 09/02/92 Dilution Factor: NA Percent Solids: NA %		
	RL	Result	Qual	RL	Result	Qual	RL	Result	Qual	RL	Result	Qual
Alkalinity - CaCO3	2.0	247	A	2.0	525	A	2.0	788	A	2.0	215	A
Ammonia - N	0.10	0.30	A	0.10	0.30	A	0.10	1.5	A	0.10	0.30	A
Biochemical Oxygen Dema	2.0	16.0	J	2.0	14	J	2.0	6.0	J	2.0	5.0	J
Chemical Oxygen Demand	5.0	98.0	A	5.0	49	A	5.0	79.0	A	5.0	39.0	A
Chloride	2.0	80.0	A	2.0	225	A	2.0	181	A	2.0	165	A
Fluoride	0.10	0.01	U	0.10	0.10	A	0.10	0.10	U	0.10	0.10	A
Hardness - CaCO3	2.0	341	A	2.0	789	A	2.0	1260	U	2.0	464	A
Hexavalent Chromium	0.10	0.01	U	0.10	0.10	U	0.10	0.10	U	0.10	0.10	U
Nitrate/Nitrite - N	0.02	0.02	U	0.02	0.05	B	0.02	0.09	B	0.02	0.50	U
Sulfate	5.0	36.0	J	5.0	114	J	5.0	332	J	5.0	33.0	J
Total Dissolved Solids	5.0	484	L	5.0	1050	L	5.0	1760	L	5.0	776	L
Total Inorganic Carbon	1.0	170	A	1.0	140	A	1.0	280	A	1.0	62.0	A
Total Kjeldahl Nitrogen	0.10	0.50	B	0.10	0.40	B	0.10	1.5	B	0.10	0.80	B
Total Organic Carbon	1.0	5.1	A	1.0	10	A	1.0	22.0	A	1.0	4.9	A
Total Phosphate - P	0.05	0.05	U	0.05	0.05	U	0.05	0.36	B	0.05	0.78	A
Total Suspended Solids	5.0	176	L	5.0	84	L	5.0	135	L	5.0	426	L
Turbidity	0.20	192	J	0.20	165	J	0.20	76.0	J	0.20	618	J

Notes: All units are mg/l except Turbidity which is NTU.
The Reporting Limits (RLs) have been adjusted for percent solids.
The Qual column indicates the qualifier applied to the result following data validation (see below).

- Qualifiers.....
- A - Acceptable (Quantitative) Data
 - J - Estimated (Semi-quantitative) Data
 - K - Biased High (Semi-quantitative) Data
 - L - Biased Low (Semi-quantitative) Data
 - B - Present in associated blanks (Semi-quantitative) Data
 - UJ - Not Detected/Estimated Data
 - U - Not Detected/Quantitative Data
 - UL - Not Detected/Biased Low
 - R - Unusable Data
 - Q - Not Applicable
 - N - Tentative Identification
 - M - Tentative Identification (result below laboratory MDL)

AR302848

October 1993

913-6773

SUMMARY OF WET CHEMISTRY CONVENTIONAL ANALYSES - PHASE 1A

BERKS COUNTY, PENNSYLVANIA

Matrix: Groundwater

Parameter	Sample Point GGG-02/1A			Sample Point GGG-03/1A			Sample Point		
	RL	Result	Qual	RL	Result	Qual	RL	Result	Qual
Alkalinity - CaCO3	2.0	104	A	2.0	64.0	A			
Ammonia - N	0.10	0.30	A	0.10	0.60	A			
Biochemical Oxygen Dema	2.0	2.0	U	2.0	2.0	U			
Chemical Oxygen Demand	5.0	7.0	B	5.0	20.0	B			
Chloride	2.0	7.0	A	2.0	4.0	A			
Fluoride	0.10	0.10	A	0.10	0.10	U			
Hardness - CaCO3	2.0	181	A	2.0	96.0	A			
Hexavalent Chromium	0.10	0.10	U	0.10	0.10	U			
Nitrate/Nitrite - N	0.02	1.84	J	0.02	0.03	B			
Sulfate	5.0	27.0	J	5.0	19.0	J			
Total Dissolved Solids	5.0	615	L	5.0	170	L			
Total Inorganic Carbon	1.0	25.0	A	1.0	24.0	A			
Total Kjeldahl Nitrogen	0.10	1.1	B	0.10	0.70	B			
Total Organic Carbon	1.0	1.0	U	1.0	1.3	A			
Total Phosphate - P	0.05	2.62	A	0.05	0.98	A			
Total Suspended Solids	5.0	412	L	5.0	950	L			
Turbidity	0.20	960	J	0.20	636	J			

Notes: All units are mg/L except Turbidity which is NTU.
 The Reporting Limits (RLs) have been adjusted for percent solids.
 The Qual column indicates the Qualifier applied to the result following data validation (see below).

- Qualifiers.....
- A - Acceptable (Quantitative) Data
 - J - Estimated (Semi-quantitative) Data
 - K - Biased High (Semi-quantitative) Data
 - L - Biased Low (Semi-quantitative) Data
 - B - Present in associated blanks (Semi-quantitative) Data
 - UJ - Not Detected/Estimated Data
 - U - Not Detected/Quantitative Data
 - UL - Not Detected/Biased Low
 - R - Unusable Data
 - Q - Not Applicable
 - M - Tentative Identification
 - M* - Tentative Identification (result below laboratory MDL)

AR302849