

NYSDEC SUPERFUND STANDBY CONTRACT
WORK ASSIGNMENT NO. D002472-14.2

**PRELIMINARY SITE ASSESSMENT REPORT
VOLUME II - SUPPORTING DOCUMENTATION**

**ENRX, INC
(FORMERLY VOELKER ANALYSIS)
CITY OF BUFFALO, NEW YORK**

SITE NO. 915150

Submitted to:

New York State Department of Environmental Conservation
Albany, New York

Submitted by:

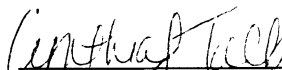
ABB Environmental Services
Portland, Maine

August 1995

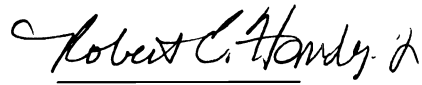
Prepared by:


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ABB Environmental
Services

Submitted by:


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Project Manager
ABB Environmental
Services

Approved by:


Robert E. Handy, Jr., P.E.
Program Manager
ABB Environmental
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ENRX, INC.
PRELIMINARY SITE ASSESSMENT REPORT
VOLUME II - SUPPORTING DOCUMENTATION

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ABB Environmental Services

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SECTION 1.0
FIELD DATA SHEETS

FIELD INSTRUMENTATION & MATERIAL QUALITY ASSURANCE RECORD

Project NYSDEC Site ENRX
 Project No. 7170-40 Sampler Signature Rick Buttl
 Date 10-3-94

Field Instrumentation Calibration Data

Equipment Type/I.D.	Battery Condition	Calibration Information
_____	_____	pH 4 _____ pH 7 _____ pH 10 _____
_____	_____	pH 4 _____ pH 7 _____ pH 10 _____
_____	_____	pH 4 _____ pH 7 _____ pH 10 _____
_____	_____	Cond. Std. _____ / _____ Cond. Std. _____ / _____ meter value
_____	_____	Cond. Std. _____ / _____ Cond. Std. _____ / _____ meter value
_____	_____	Cond. Std. _____ / _____ Cond. Std. _____ / _____ meter value
Dissolved Oxygen		
_____	_____	Avg. Winkler Value _____ ppm Meter Value _____ ppm
Redox		
_____	_____	Zobell Sol. Value _____ Meter Value _____
Photoionization Meter		
<u>Thermo Env. 580B</u>	<u>OK</u>	Zero/Zero Air? ☒ Yes ☐ No Span Gas Value <u>100</u> ppm Equiv.
<u>(HAZCO # 35114)</u>		Meter Value <u>100</u> ppm Equiv.
_____	_____	Zero/Zero Air? ☐ Yes ☐ No Span Gas Value _____ ppm Equiv.
		Meter Value _____ ppm Equiv.
Other		
<u>HAZCO # 1545 MX-241 (EL) O₂</u>	<u>OK</u>	<u>Precalibrated. Amb = 20.8% O₂</u>

Fluids/Materials Record

Deionized Water Source: _____ ECJ Staging Portable System Other
 Trip Blank Water Source: _____ ECJ Lab; Lot No. _____
 _____ Other; Type _____ ID _____
 Decontamination Fluids: _____ Methyl Hydrate; Lot No. _____
 _____ Other; Type _____ ID _____
 HNO₃/DI Rinse Solution: _____ ECJ Staging; Lot No. _____
 Filtration Paper ID: (In Line) Manuf/Type _____ Lot No. _____ / _____
 (Vacuum) Manuf/Type _____ Lot No. _____ / _____
 Chemicals Used: HNO₃ Lot No. _____ ZnAOC Lot No. _____
 H₂SO₄ Lot No. _____ Other Lot No. _____
 HCL Lot No. _____ Other Lot No. _____
 NaOH Lot No. _____

FIELD INSTRUMENTATION & MATERIAL QUALITY ASSURANCE RECORD

Project NYSDDEC Site ENRX
 Project No. 7170-40 Sampler Signature David Wilkie
 Date 10-4-94

Field Instrumentation Calibration Data

Equipment Type/I.D.	Battery Condition	Calibration Information
(HAZCO 3896) <u>HORIBA WATER CHECKER U-10</u> <u>(mF 304001)</u>	<u>GOOD</u>	pH 4 ☒ pH 7 ☐ pH 10 ☐ pH 4 ☐ pH 7 ☐ pH 10 ☐ pH 4 ☐ pH 7 ☐ pH 10 ☐ Cond. Std. ☐ / ☐ Cond. Std. ☐ / ☐ Cond. Std. ☐ / ☐ Cond. Std. ☐ / ☐ Cond. Std. ☐ / ☐ Cond. Std. ☐ / ☐
Dissolved Oxygen/Salinity/turbid/Temp <u>HORIBA U-10</u>	<u>GOOD</u>	Avg. Winkler Value ppm Meter Value <u>5.86</u> ppm Zobell Sol. Value Meter Value
Redox		
Photoionization Meter <u>THERM ENVIR. (HAZCO #3914)</u> <u>580 P</u>		Zero/Zero Air? ☒ Yes ☐ No Span Gas Value <u>100</u> ppm Equiv. } <u>see HAZCO</u> Meter Value <u>100</u> ppm Equiv. } <u>Cal sheet.</u> Zero/Zero Air? ☐ Yes ☐ No Span Gas Value ppm Equiv. Meter Value ppm Equiv.
Other <u>Brüel + Kjaer Møllgas Monitor</u> <u>HAZCO MX-241 LEL/O2</u>	<u>OK</u>	<u>Pre-Calibrated</u> <u>(#1565- see HAZCO Cal. sheet)</u>

Fluids/Materials Record

Deionized Water Source: ☒ ECJ Staging Portable System Other A/B/B SOURCE
 Trip Blank Water Source: ☐ ECJ Lab; Lot No.
☒ Other; Type NYTEST ENV. ID ENQT001XXX94XX
 Decontamination Fluids: ☐ Methyl Hydrate; Lot No.
☒ Other; Type LIQUINOX ID
 HNO₃/DI Rinse Solution: ☐ ECJ Staging; Lot No.
 Filtration Paper ID: (In Line) Manuf/Type Lot No.
 (Vacuum) Manuf/Type Lot No.
 Chemicals Used: HNO₃ Lot No. ZnAOC Lot No.
 H₂SO₄ Lot No. Other Lot No.
 HCL Lot No. Other Lot No.
 NaOH Lot No.

E.C. JORDAN, CO.

FIELD INSTRUMENTATION & MATERIAL QUALITY ASSURANCE RECORD

Project NYSDEC Site ENRX
 Project No. 7170-40 Sampler Signature BE Buttl
 Date 10-5-94

Field Instrumentation Calibration Data

Equipment Type/I.D.

Battery Condition

Calibration Information

~~Auto~~ calibration Fluid used

Horiba U-10 W.Q. checker
(HAZCO 3896)

OK

pH 4 ✓ pH 7 pH 10

pH 4 pH 7 pH 10

pH 4 pH 7 pH 10

Cond. Std. NO std. Cond. Std. NO std. meter value

Cond. Std. / Cond. Std. / meter value

Cond. Std. / Cond. Std. / meter value

Dissolved Oxygen / salinity / turbidity / Temp

Horiba U-10

OK

Avg. Winkler Value ppm Meter Value 5.86 ppm

Redox

Zobell Sol. Value Meter Value

Photoionization Meter

Thermo Environmental 580B
(HAZCO # 3914)

OK

Zero/Zero Air? ☒ Yes ☐ No Span Gas Value 100 ppm Equiv.

Meter Value 100 ppm Equiv.

Zero/Zero Air? ☐ Yes ☐ No Span Gas Value ppm Equiv.

Meter Value ppm Equiv.

Other

Bauer & Taer MultiGas Mon.
HAZCO MX-241 LEL/O2

OK

OK

Pre-calibrated

(#1565 - See HAZCO sheet)

Fluids/Materials Record

Deionized Water Source: ☒ ECJ Staging Portable System Other

Trip Blank Water Source: ECJ Lab; Lot No.

☒ Other; Type NYTEST Env.

ID As yesterday

Decontamination Fluids: Methyl Hydrate; Lot No.

☒ Other; Type Liquinox

ID

HNO₃/DI Rinse Solution: ECJ Staging; Lot No.

Filtration Paper ID: (In Line) Manuf/Type Lot No. /

(Vacuum) Manuf/Type Lot No. /

Chemicals Used: HNO₃ Lot No. Lab Pre-preserved ZnAOC Lot No.

H₂SO₄ Lot No. Other Lot No.

HCL Lot No. Other Lot No.

NaOH Lot No. Lab Pre-preserved

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1302 Filter and Calibration Data

CUSTOMER: ABB Environmental

SERIAL NUMBER: 1617171

Filter Number: UA 0936

Filter Position: A

Filter Bank #: 1

Zero Calibration:

Date: 09/27/94

Calibration Gas: Zero air

Humidity Interference Calibration:

Date: 09/27/94

Calibration Gas: Humidified zero air

Dewpoint (°C): 13.8

Span Calibration:

Date: 00/00/94

Calibration Gas:

Concentration (ppm):

Source:

Ambient Temp. (°C):

Ambient Pressure (mmHg):

Measurment Results:

Input Concentration(ppm)

1302 Measurement Reading(ppm)

1302 Filter and Calibration Data

CUSTOMER: ABB Environmental

SERIAL NUMBER: 1617171

Filter Number: UA 0970

Filter Position: B

Filter Bank #: 1

Zero Calibration:

Date: 09/27/94

Calibration Gas: Zero air

Humidity Interference Calibration:

Date: 09/27/94

Calibration Gas: Humidified zero air

Dewpoint (°C): 13.8

Span Calibration:

Date: 09/28/94

Calibration Gas: *trans*-1,2-Dichloroethylene

Concentration (ppm): 265

Source: Liquid injection into a Tedlar bag

Ambient Temp. (°C): 22.0

Ambient Pressure (mmHg): 738

Measurment Results:

<u>Input Concentration(ppm)</u>	<u>1302 Measurement Reading(ppm)</u>
265	263 ± 1
120	123 ± 1
48.2	49.1 ± 0.1

1302 Filter and Calibration Data

CUSTOMER: ABB Environmental

SERIAL NUMBER: 1617171

Filter Number: UA 0972

Filter Position: C

Filter Bank #: 1

Zero Calibration:

Date: 09/27/94

Calibration Gas: Zero air

Humidity Interference Calibration:

Date: 09/27/94

Calibration Gas: Humidified zero air

Dewpoint (°C): 13.8

Span Calibration:

Date: 09/29/94

Calibration Gas: 1,1,1-Trichloroethane

Concentration (ppm): 243

Source: Liquid injection into a Tedlar bag

Ambient Temp. (°C): 22.0

Ambient Pressure (mmHg): 738

Measurment Results:

<u>Input Concentration(ppm)</u>	<u>1302 Measurement Reading(ppm)</u>
243	244 ± 1
131	128 ± 1
37.3	37.5 ± 0.2

1302 Filter and Calibration Data

CUSTOMER: ABB Environmental

SERIAL NUMBER: 1617171

Filter Number: UA 0974

Filter Position: D

Filter Bank #: 1

Zero Calibration:

Date: 09/27/94

Calibration Gas: Zero air

Humidity Interference Calibration:

Date: 09/27/94

Calibration Gas: Humidified zero air

Dewpoint (°C): 13.8

Span Calibration:

Date: 09/29/94

Calibration Gas: 1,1-Dichloroethylene

Concentration (ppm): 256

Source: Liquid injection into a Tedlar bag

Ambient Temp. (°C): 22.0

Ambient Pressure (mmHg): 738

Measurment Results:

<u>Input Concentration(ppm)</u>	<u>1302 Measurement Reading(ppm)</u>
256	255 ± 1
116	120 ± 1
46.6	47.9 ± 0.1

1302 Filter and Calibration Data

CUSTOMER: ABB Environmental

SERIAL NUMBER: 1617171

Filter Number: UA 0980

Filter Position: E

Filter Bank #: 1

Zero Calibration:

Date: 09/27/94

Calibration Gas: Zero air

Humidity Interference Calibration:

Date: 09/27/94

Calibration Gas: Humidified zero air

Dewpoint (°C): 13.8

Span Calibration:

Date: 09/30/94

Calibration Gas: Carbon tetrachloride

Concentration (ppm): 154

Source: Liquid injection into a Tedlar bag

Ambient Temp. (°C): 22.0

Ambient Pressure (mmHg): 740

Measurment Results:

<u>Input Concentration(ppm)</u>	<u>1302 Measurement Reading(ppm)</u>
154	154 ± 1
77.1	77.8± 0.1
38.6	38.8 ± 0.1

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HAZCO Services, Inc.

Instrument Description

580B

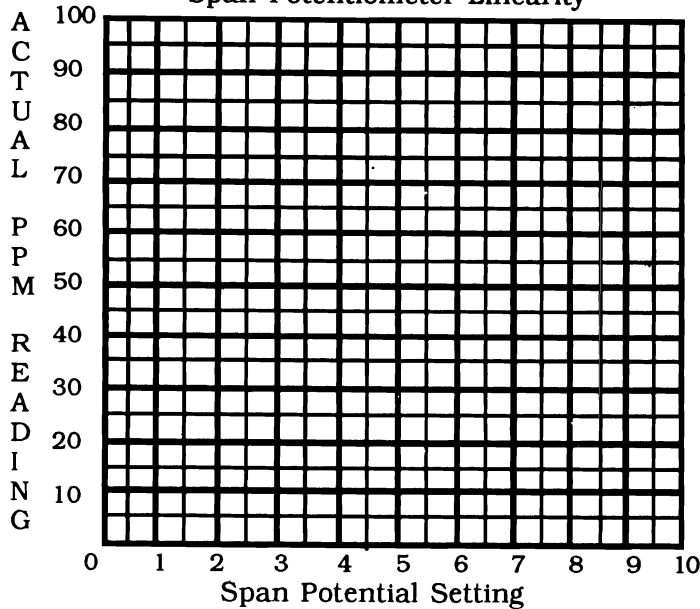
Mfg. Serial #: 40105-262

Calibration Date: 9-30-94

HAZCO Serial #: 3914

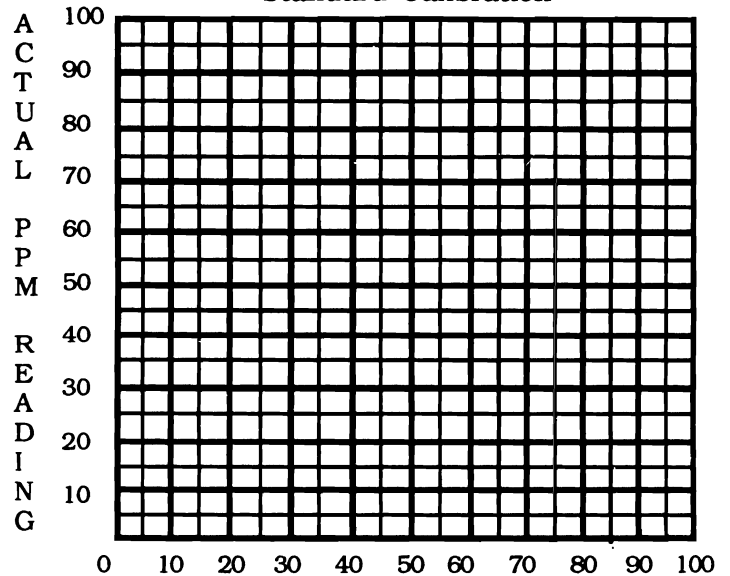
Technician: CDG

Span Potentiometer Linearity



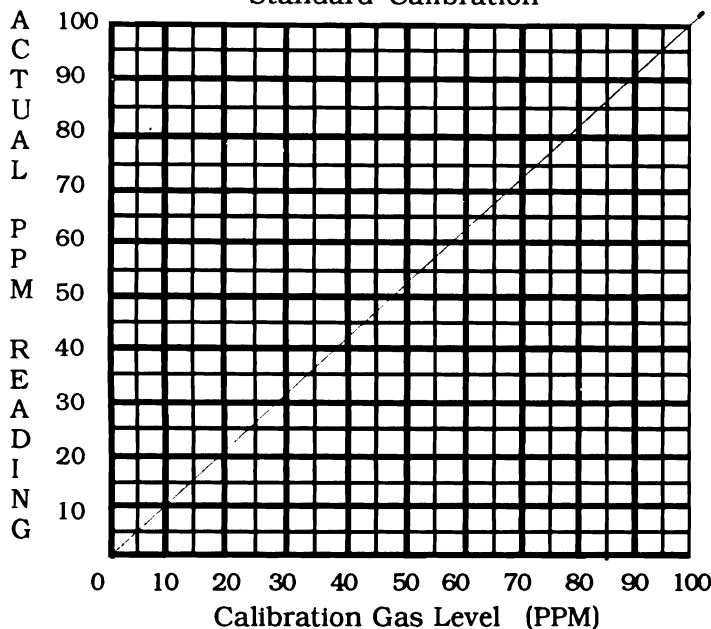
Calibration Points

Standard Calibration



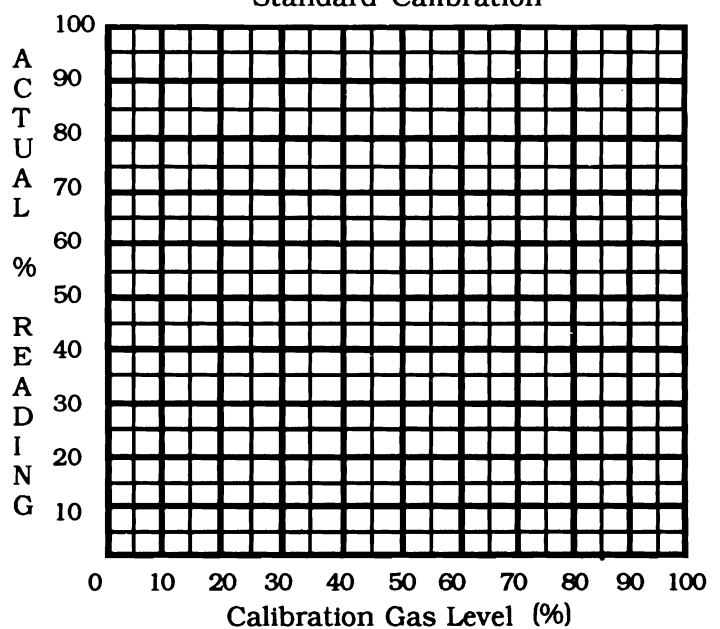
Calibration Points

Standard Calibration



Calibration Points 00, 104-104

Standard Calibration



Calibration Points

Please Return Equipment to

- ☒ HAZCO Services, Inc., 2006 Springboro West, Dayton, OH 45439, 513-293-2700
- ☐ HAZCO Services, Inc., 2550 Eisenhower Ave., Valley Forge, PA 19481, 610-650-8900
- ☐ HAZCO Services, Inc., 15353 East Hinsdale Circle, Suite D, Englewood, CO 80112, 303-766-1501
- ☐ HAZCO Services, Inc., 1735 East Wilshire Blvd., Suite 802, Santa Ana, CA 92705, 714-558-1475
- ☐ HAZCO Services, Inc., 50 Campus Plaza, Edison, NJ 08837-3911, 908-417-0660,
- ☐ HAZCO Services, Inc., 115-A Flint Road, Oak Ridge, TN 37830, 615-482-3940



HAZCO Services, Inc.

Instrument Description

MX241

Mfg. Serial #:

Calibration Date:

9-30-94

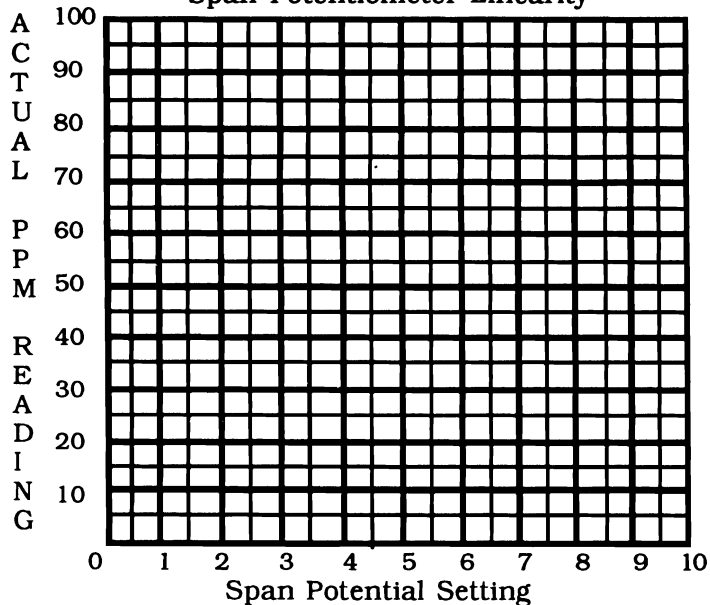
HAZCO Serial #:

1565

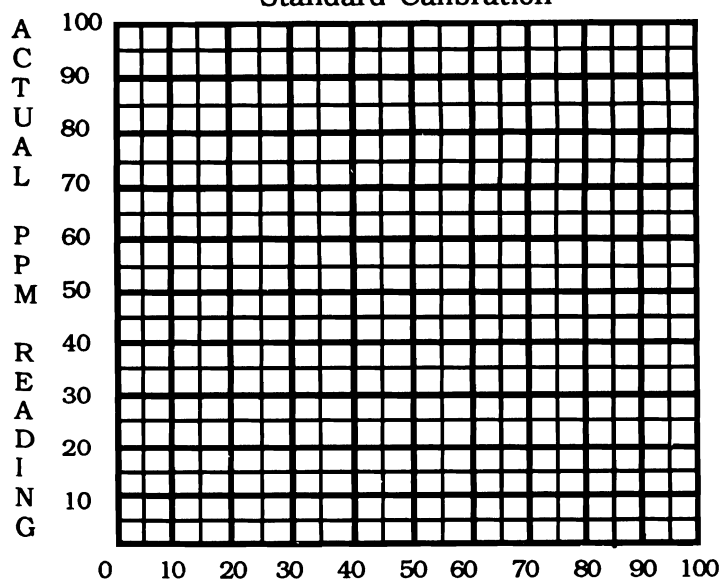
Technician:

SRR

Span Potentiometer Linearity



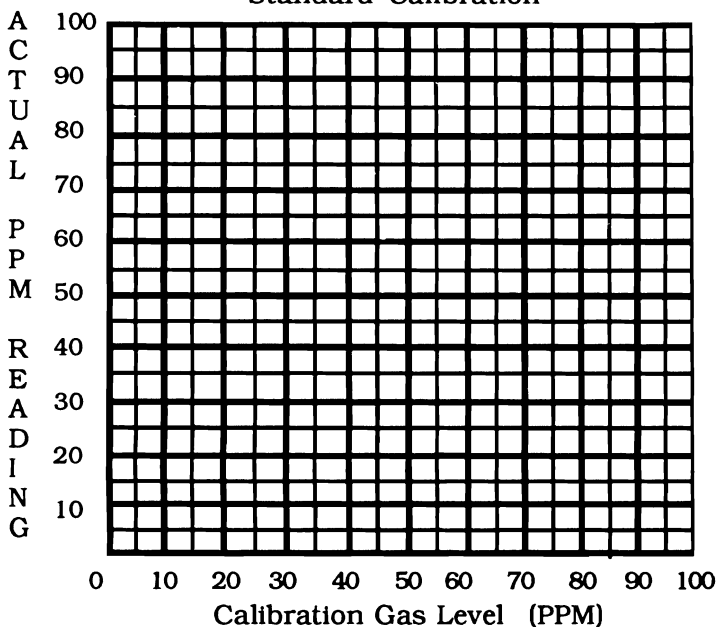
Standard Calibration



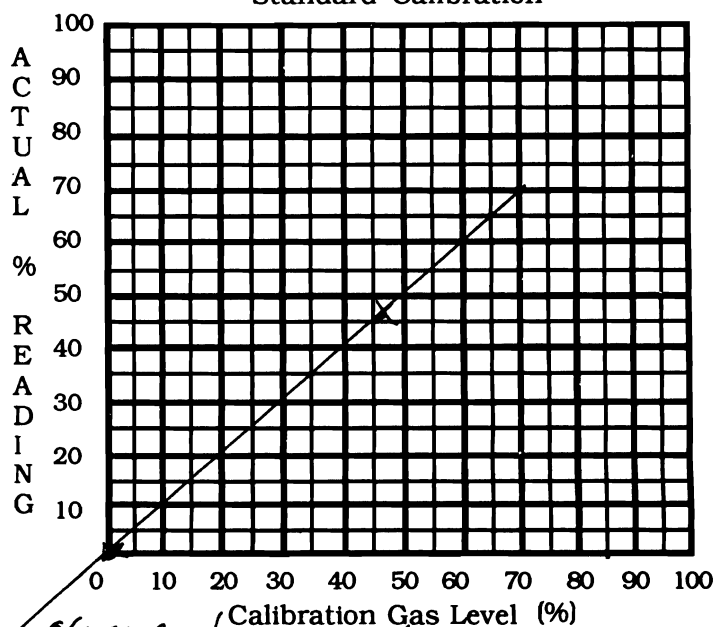
Calibration Points

Calibration Points

Standard Calibration



Standard Calibration



Calibration Points

%LEL Pentane
Calibration Points 0-0 / 46-46

Please Return Equipment to

- ☒ HAZCO Services, Inc., 2006 Springboro West, Dayton, OH 45439, 513-293-2700
- ☐ HAZCO Services, Inc., 2550 Eisenhower Ave., Valley Forge, PA 19481, 610-650-8900
- ☐ HAZCO Services, Inc., 15353 East Hinsdale Circle, Suite D, Englewood, CO 80112, 303-766-1501
- ☐ HAZCO Services, Inc., 1735 East Wilshire Blvd., Suite 802, Santa Ana, CA 92705, 714-558-1475
- ☐ HAZCO Services, Inc., 50 Campus Plaza, Edison, NJ 08837-3911, 908-417-0660,
- ☐ HAZCO Services, Inc., 115-A Flint Road, Oak Ridge, TN 37830, 615-482-3940

FIELD INSTRUMENTATION & MATERIAL QUALITY ASSURANCE RECORD

Project ENRY INC. Site ENRY INC.
 Project No. 7170-40 Sampler Signature Jim D. Gyl
 Date 10-27-94

Field Instrumentation Calibration Data

Equipment Type/I.D.	Battery Condition	Calibration Information
_____	_____	pH 4 _____ pH 7 _____ pH 10 _____
_____	_____	pH 4 _____ pH 7 _____ pH 10 _____
_____	_____	pH 4 _____ pH 7 _____ pH 10 _____
_____	_____	Cond. Std. _____/_____ Cond. Std. _____/_____
_____	_____	Cond. Std. _____/_____ Cond. Std. _____/_____
_____	_____	Cond. Std. _____/_____ Cond. Std. _____/_____
Dissolved Oxygen	_____	Avg. Winkler Value _____ ppm Meter Value _____ ppm
Redox	_____	Zobell Sol. Value _____ Meter Value _____
Photoionization Meter	_____	Zero/Zero Air? ☐ Yes ☒ No Span Gas Value <u>100</u> ppm Equiv.
<u>TE 580B, OVM #5315 OK</u>		Meter Value <u>104.5</u> ppm Equiv.
<u>(Rental From HAZCO)</u>		Zero/Zero Air? ☐ Yes ☐ No Span Gas Value _____ ppm Equiv.
		Meter Value _____ ppm Equiv.
Other		
<u>DUST Monitors</u>	<u>OK</u>	<u>RENTED FROM HAZCO</u>
<u>MINIRAM</u>		

Fluids/Materials Record

Deionized Water Source: ☒ ECJ Staging ☐ Portable System ☐ Other

Trip Blank Water Source: _____ ECJ Lab; Lot No. _____

_____ Other; Type _____ ID _____

Decontamination Fluids: _____ Methyl Hydrate; Lot No. _____

☒ Other; Type POTABLE WATER & LIQUINOX

HNO₃/DI Rinse Solution: _____ ECJ Staging; Lot No. _____

Filtration Paper ID: (In Line) Manuf/Type _____ Lot No. _____/_____

(Vacuum) Manuf/Type _____ Lot No. _____/_____

Chemicals Used: HNO₃ Lot No. * ☒ LAB SUPPLIED ZnAOC Lot No. _____

H₂SO₄ Lot No. _____ Other Lot No. _____

HCL Lot No. * ☒ LAB SUPPLIED Other Lot No. _____

NaOH Lot No. * ☒ LAB SUPPLIED

* = LAB BOTTLES PRE-PRESERVED FROM LYTIST

E.C. JORDAN, CO.

FIELD INSTRUMENTATION & MATERIAL QUALITY ASSURANCE RECORD

Project _____ Site ENRX, INC.
 Project No. 7170-40 Sampler Signature John S. Zyl
 Date 10-28-94

Field Instrumentation Calibration Data

Equipment Type/I.D.	Battery Condition	Calibration Information
_____	_____	pH 4 _____ pH 7 _____ pH 10 _____
_____	_____	pH 4 _____ pH 7 _____ pH 10 _____
_____	_____	pH 4 _____ pH 7 _____ pH 10 _____
_____	_____	Cond. Std. _____ / _____ Cond. Std. _____ / _____ meter value
_____	_____	Cond. Std. _____ / _____ Cond. Std. _____ / _____ meter value
_____	_____	Cond. Std. _____ / _____ Cond. Std. _____ / _____ meter value
Dissolved Oxygen		Avg. Winkler Value _____ ppm Meter Value _____ ppm
Redox		Zobell Sol. Value _____ Meter Value _____
Photoionization Meter		Zero/Zero Air? ☐ Yes ☐ No Span Gas Value _____ ppm Equiv.
<u>580 B OVA</u>	<u>OK</u>	Meter Value _____ ppm Equiv.
		Zero/Zero Air? ☒ Yes ☐ No Span Gas Value <u>100</u> ppm Equiv.
		Meter Value <u>99.9</u> ppm Equiv.
Other		<u>(FLUCTUATED)</u>

Fluids/Materials Record

Deionized Water Source: ☒ ECJ Staging Portable System Other
 Trip Blank Water Source: ☒ ECJ Lab; Lot No. _____
 Other; Type LAB SUPPLIED ID _____
 Decontamination Fluids: ☒ Methyl Hydrate; Lot No. _____
 Other; Type DI, POTABLE H2O ID LIQUOR
 HNO₃/DI Rinse Solution: _____ ECJ Staging; Lot No. _____
 Filtration Paper ID: (In Line) Manuf/Type _____ Lot No. _____ / _____
 (Vacuum) Manuf/Type _____ Lot No. _____ / _____
 Chemicals Used: HNO₃ Lot No. _____ ZnAOC Lot No. _____
 H₂SO₄ Lot No. _____ Other Lot No. _____
 HCL Lot No. _____ Other Lot No. _____
 NaOH Lot No. _____

ALL LAB SUPPLIED W/ SAMPLE BOTTLES

FIELD INSTRUMENTATION & MATERIAL QUALITY ASSURANCE RECORD

Project _____ Site ENRY, INC.
 Project No. 7170-40 Sampler Signature John A. Lyly
 Date 10-31-94

Field Instrumentation Calibration Data

Equipment Type/I.D.	Battery Condition	Calibration Information
<u>Horiba Water Checker</u>	<u>OK</u>	<u>AUTO CALIBRATION</u>
<u>U-10 #5090 HAZCO</u>		<u>FEATURE W/</u>
		<u>INSTRUMENT</u>
		pH 4 _____ pH 7 _____ pH 10 _____
		pH 4 _____ pH 7 _____ pH 10 _____
		pH 4 _____ pH 7 _____ pH 10 _____
		Cond. Std. _____ / _____ meter value
		Cond. Std. _____ / _____ meter value
		Cond. Std. _____ / _____ meter value
Dissolved Oxygen		Avg. Winkler Value _____ ppm Meter Value _____ ppm
Redox		Zobell Sol. Value _____ Meter Value _____
Photoionization Meter		Zero/Zero Air? ☐ Yes ☐ No Span Gas Value _____ ppm Equiv.
		Meter Value _____ ppm Equiv.
<u>TE 580B OVM</u>	<u>OK</u>	<u>AMBIENT</u>
<u>#5315 HAZCO</u>		<u>Zero/Zero Air?</u> ☒ Yes ☐ No Span Gas Value <u>100</u> ppm Equiv.
		Meter Value <u>105 to 99</u> ppm Equiv.
Other		<u>FLUCTUATED</u>

Fluids/Materials Record

Deionized Water Source: ☒ ECJ Staging ☐ Portable System ☐ Other _____
 Trip Blank Water Source: _____ ECJ Lab; Lot No. _____
 _____ Other; Type _____ ID _____
 Decontamination Fluids: ☒ Methyl Hydrate; Lot No. _____
☒ Other; Type DI, POTABLE H₂O to LIQUINOX
 HNO₃/DI Rinse Solution: _____ ECJ Staging; Lot No. _____
 Filtration Paper ID: (In Line) Manuf/Type _____ Lot No. _____ / _____
 (Vacuum) Manuf/Type _____ Lot No. _____ / _____
 Chemicals Used: HNO₃ Lot No. _____ ZnAOC Lot No. _____
 H₂SO₄ Lot No. _____ Other Lot No. _____
 HCL Lot No. _____ Other Lot No. _____
 NaOH Lot No. _____

FIELD INSTRUMENTATION & MATERIAL QUALITY ASSURANCE RECORD

Project _____ Site ENRX, INC.
 Project No. 7170-40 Sampler Signature John W. Lyby
 Date 11-1-94

Field Instrumentation Calibration Data

Equipment Type/I.D.

Battery
Condition

Calibration Information

Very Heavy Rain Today

Hanna Water Checker
U-10, #5090 HAZCO

OK

pH 4 _____ pH 7 _____ pH 10 _____

pH 4 _____ pH 7 _____ pH 10 _____

pH 4 _____ pH 7 _____ pH 10 _____

Cond. Std. _____ / _____ Cond. Std. _____ / _____ meter value

Cond. Std. _____ / _____ Cond. Std. _____ / _____ meter value

Cond. Std. _____ / _____ Cond. Std. _____ / _____ meter value

*AUTO CALIBRATION
PERFORMED*

Dissolved Oxygen

Avg. Winkler Value _____ ppm Meter Value _____ ppm

Redox

Zobell Sol. Value _____ Meter Value _____

Photoionization Meter

Zero/Zero Air? ☐ Yes ☐ No Span Gas Value _____ ppm Equiv.

Meter Value _____ ppm Equiv.

Zero/Zero Air? ☐ Yes ☐ No Span Gas Value _____ ppm Equiv.

Meter Value _____ ppm Equiv.

Other

Fluids/Materials Record

Deionized Water Source: ☒ ECJ Staging ☐ Portable System ☐ Other

Trip Blank Water Source: _____ ECJ Lab; Lot No. _____

Other; Type _____ ID _____

Decontamination Fluids: ☒ Methyl Hydrate; Lot No. _____

Other; Type DI, POTABLE H₂O LIQUINOX

HNO₃/DI Rinse Solution: _____ ECJ Staging; Lot No. _____

Filtration Paper ID: (In Line) Manuf/Type _____ Lot No. _____ / _____

(Vacuum) Manuf/Type _____ Lot No. _____ / _____

Chemicals Used: HNO₃ Lot No. _____ ZnAOC Lot No. _____

H₂SO₄ Lot No. _____ Other Lot No. _____

HCL Lot No. _____ Other Lot No. _____

NaOH Lot No. _____

ABB Environmental Services

FIELD INSTRUMENTATION & MATERIAL QUALITY ASSURANCE RECORD

Project ENRX, Inc. Site ENRX Inc
 Project No. 7170-40 Sampler Signature Rick Buttl
 Date 11-30-94

Field Instrumentation Calibration Data

Equipment Type/I.D.	Battery Condition	Calibration Information
<u>Horiba U-10 WQ checker</u>	<u>OK</u>	pH 4 <u>✓</u> pH 7 <u> </u> pH 10 <u> </u>
<u>(Hazeo #)</u>	<u> </u>	pH 4 <u> </u> pH 7 <u> </u> pH 10 <u> </u>
<u>*Autocal. Sol'n*</u>	<u> </u>	pH 4 <u> </u> pH 7 <u> </u> pH 10 <u> </u>
<u> </u>	<u> </u>	Cond. Std. <u>Not Avail.</u> Cond. Std. <u>Not Avail.</u> meter value
<u> </u>	<u> </u>	Cond. Std. <u> </u> / <u> </u> Cond. Std. <u> </u> / <u> </u> meter value
<u> </u>	<u> </u>	Cond. Std. <u> </u> / <u> </u> Cond. Std. <u> </u> / <u> </u> meter value
Dissolved Oxygen /		
<u>Horiba - U 10 (Hazeo #)</u>	<u>OK</u>	Avg. Winkler Value <u> </u> ppm Meter Value <u> </u> ppm
Redox		
<u> </u>	<u> </u>	Zobell Sol. Value <u> </u> Meter Value <u> </u>
Photoionization Meter		
<u>Thermo Environmental</u>	<u> </u>	Zero/Zero Air? ☐ Yes ☐ No Span Gas Value <u> </u> ppm Equiv. Meter Value <u> </u> ppm Equiv.
<u> </u>	<u> </u>	Zero/Zero Air? ☐ Yes ☐ No Span Gas Value <u> </u> ppm Equiv. Meter Value <u> </u> ppm Equiv.
Other		
<u> </u>	<u> </u>	<u> </u>

Fluids/Materials Record

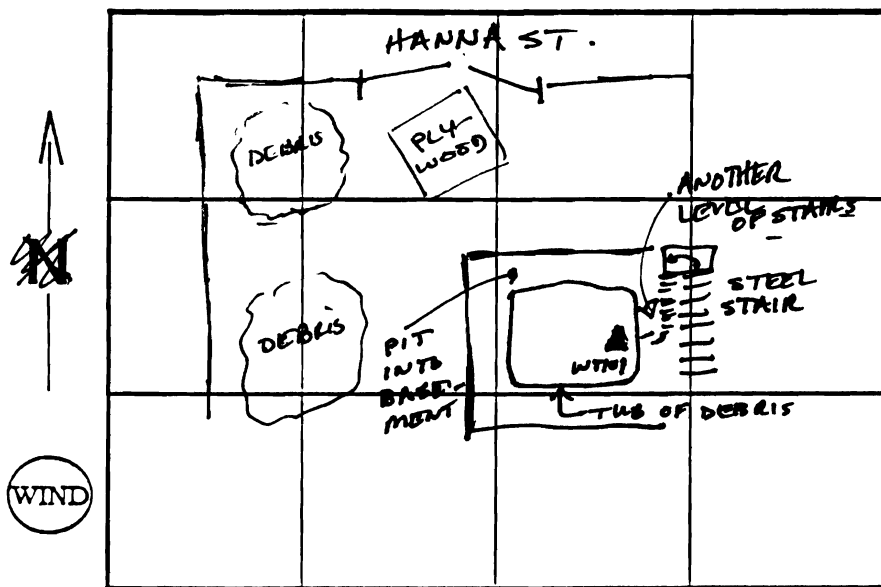
Deionized Water Source: ☒ ECJ Staging Portable System Other
 Trip Blank Water Source: ECJ Lab; Lot No.
☒ Other; Type NYTEST ID ENGT002XXX94XX
 Decontamination Fluids: Methyl Hydrate; Lot No.
☒ Other; Type Liquinox ID
 HNO₃ /DI Rinse Solution: ECJ Staging; Lot No.
 Filtration Paper ID: (In Line) Manuf/Type Lot No. /
 (Vacuum) Manuf/Type Lot No. /
 Chemicals Used: HNO₃ Lot No. Lab prepreserved ZnAOC Lot No.
 H₂SO₄ Lot No. Other Lot No.
 HCL Lot No. Other Lot No.
 NaOH Lot No. Lab prepreserved

SUMP/DRY WELL/STRUCTURE SAMPLING RECORD

PAGE 1 OF 1

SITE ENRX STRUCTURE TYPE Still bottoms Tank
 STRUCTURE ID WT 101 DATE 10-4-94 TIME 1640 END 1700
 COORDINATES _____

PLAN VIEW OF STRUCTURE SITE W/ DIMENSIONS



SCALE 1 INCH= NTS FEET

CREW MEMBERS:

1. D. LOVELLOY
2. C. TALBOT
3. D. WILKIE
4. B. BUTLER
- 5.
- 6.

MONITORING EQUIPMENT:

PI Meter ☒ N
 Explosive Gas ☒ N
 Avail. Oxygen ☒ N
 OVA Y N
 Other N/A

Photographs, Roll N/A

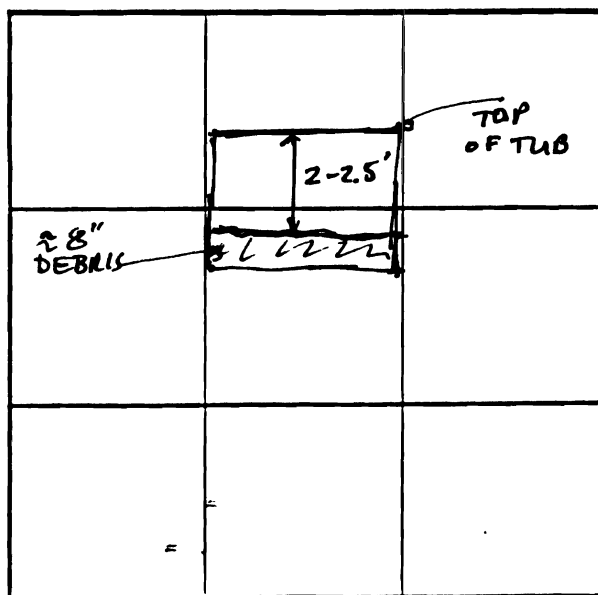
Exposure N/A

HEALTH AND SAFETY

Protection Level B
 Initial PI 0. ppm
 Initial LEL — %
 Initial O₂ 20.6 %

PID Sample = 0 ppm

CROSS SECTION OF STRUCTURE



SCALE 1 INCH= NTS FEET

Logged by D Wilkie

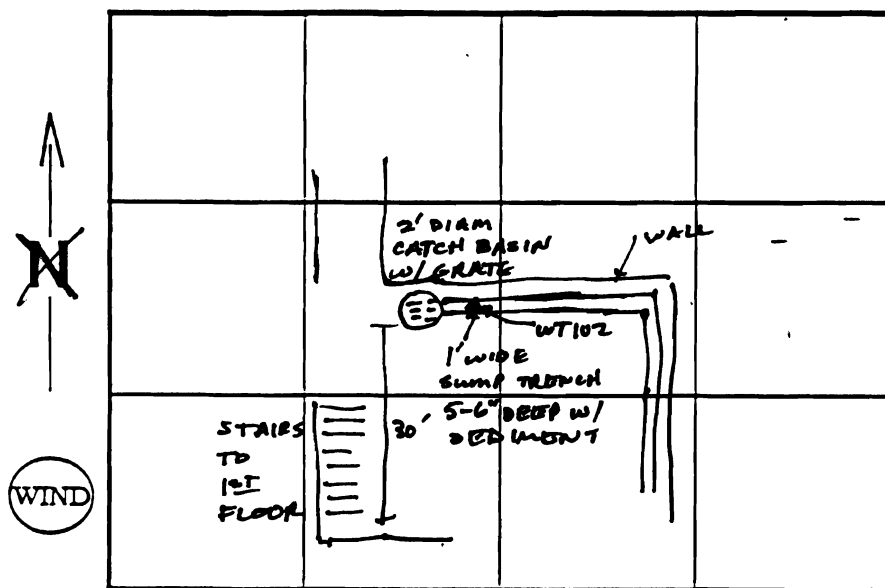
Checked by B Butler

SUMP/DRY WELL/STRUCTURE SAMPLING RECORD

PAGE 1 OF 1

SITE ENRX, Inc. STRUCTURE TYPE sump.
 STRUCTURE ID WT102 DATE 10-4-94 TIME 1605 END 1645
 COORDINATES _____

PLAN VIEW OF STRUCTURE SITE W/ DIMENSIONS



SCALE 1 INCH = NTS FEET

CREW MEMBERS:

1. D. LOVEJOY
2. C. TALBOT
3. D. WILKIE
4. B. BUTLER
- 5.
- 6.

MONITORING EQUIPMENT:

PI Meter Y N
 Explosive Gas Y N
 Avail. Oxygen Y N
 OVA Y N
 Other _____

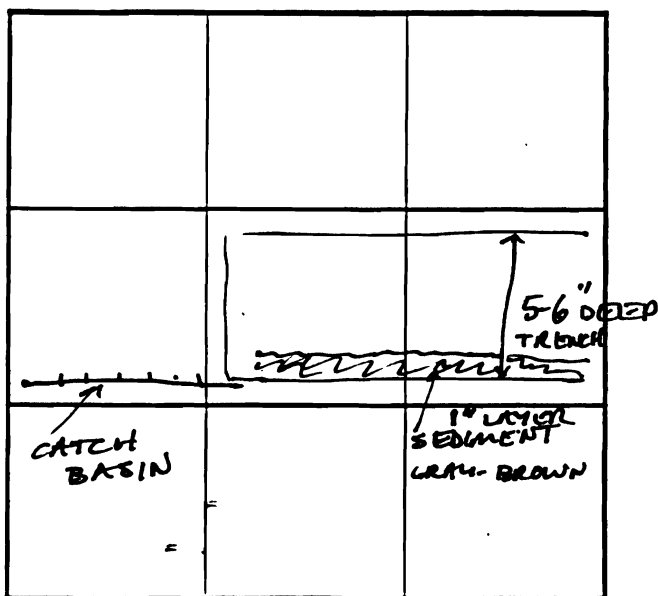
Photographs. Roll _____

Exposure _____

HEALTH AND SAFETY

Protection Level B
 Initial PI 25-30 ppm
 Initial LEL 1 %
 Initial O₂ 20.6 %

CROSS SECTION OF STRUCTURE



SCALE 1 INCH = NTS FEET

Logged by D. Wilkie

Checked by B. Butler

SURFACE SOIL SAMPLE DATA RECORD

Project: ENRX Site: First Floor
 Project Number: 7170-40 Date: 10-4-74
 Sample Location ID: WT-103/104
 Time: Start: 1700 End: 1730 Signature of Sampler: D. Wilkins

SOIL SAMPLE

DEPTH OF SAMPLE 0-1"

EQUIPMENT USED FOR COLLECTION:

☐ HAND AUGER
☐ S.S. SPLT SPOON
☐ SHOVEL
☒ HAND SPOON
☐ ALUMINUM PANS
☐ SS BUCKET

DECONTAMINATION FLUIDS USED:

☒ ALL USED
☐ ETHYL ALCOHOL
☐ 25% METHANOL/ 75% ASTM TYPE II WATER
☒ DEIONIZED WATER
☒ LIQUINOX SOLUTION
☐ HEXANE
☐ HNO₃ SOLUTION
☐ POTABLE WATER
☐ NONE

TYPE OF SAMPLE COLLECTED:

☒ DISCRETE
☐ COMPOSITE

SOIL TYPE:

☐ CLAY
☐ SAND
☐ ORGANIC
☐ GRAVEL

SAMPLE OBSERVATIONS:

☐ ODOR dark gray w/
☐ COLOR some white spots
☐ most

☒ other: sludge

FIELD GC DATA: ☐ FIELD DUPLICATE COLLECTED
 DUPLICATE ID _____

SAMPLE LOCATION SKETCH:

☒ YES see below
☐ NO

SAMPLES COLLECTED

✓ IF REQUIRED AT THIS LOCATION	MATRIX		✓ IF PRESERVED WITH ACID-BASE	VOLUME REQUIRED	✓ IF SAMPLE COLLECTED	SAMPLE BOTTLE IDS
	SURFACE WATER	SEDIMENT				
☒ TCL UCC	☐	☐	☐	_____	☒	_____
☒ TCL SWCC	☐	☐	☐	_____	☒	_____
☒ TCL Pest/Res	☐	☐	☐	_____	☒	_____
☒ TCL Inorg.	☐	☐	☐	_____	☒	_____
☒ EPTOX metab	☐	☐	☐	_____	☒	_____
☒ Inorganic	☐	☐	☐	_____	☒	_____

NOTES/SKETCH

HANNA ST
 Sample location
 WT-103
 Pit to
 Basement
 holes in ceiling
 (STILL debris)
 WT-104
 104

SURFACE SOIL SAMPLE DATA RECORD

Project: ENRX, INC.
 Project Number: 770-40
 Sample Location ID: WT-105
 Time: Start: 0845 End: 0955

Site: Second Floor
 Date: 10-5-94
 Signature of Sampler: BK Buttl for Wilke

SOIL SAMPLE

DEPTH OF SAMPLE 1" PILE ON

FLOOR
- CAULKED ON
TO FLOOR,
NEEDED
TO BE SCRAPED
OFF FLOOR.

EQUIPMENT USED FOR COLLECTION:

- ☐ HAND AUGER
- ☐ S.S. SPLIT SPOON
- ☐ SHOVEL
- ☒ HAND SPOON
- ☐ ALUMINUM PANS
- ☐ SS BUCKET

TYPE OF SAMPLE COLLECTED:

- ☐ DISCRETE
- ☐ COMPOSITE

SAMPLE OBSERVATIONS:

- ☐ ODOR
- ☐ COLOR DARK BROWN
- ☐ LIGHT MATERIAL
- POSSIBLE LIME OR SODA ASH

DECONTAMINATION FLUIDS USED:

- ☒ ALL USED
- ☐ ETHYL ALCOHOL
- ☐ 25% METHANOL/ 75% ASTM TYPE II WATER
- ☐ DEIONIZED WATER
- ☒ LIQUINOX SOLUTION
- ☐ HEXANE
- ☐ HNO₃ SOLUTION
- ☐ POTABLE WATER
- ☐ NONE

SOIL TYPE:

- ☐ CLAY
- ☒ SAND - POWDER
- ☐ ORGANIC
- ☐ GRAVEL

FIELD GC DATA: ☐ FIELD DUPLICATE COLLECTED
 DUPLICATE ID

SAMPLE LOCATION SKETCH:

- ☐ YES
- ☐ NO

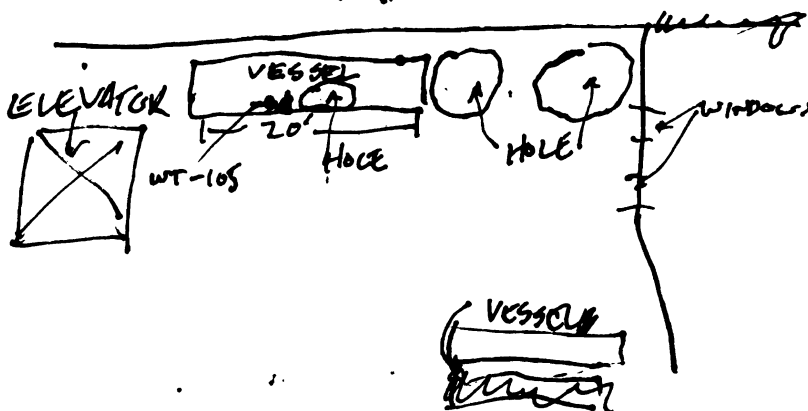
P10=0.

SAMPLES COLLECTED

/ IF REQUIRED AT THIS LOCATION	MATRIX		/ IF PRESERVED WITH ACID-BASE	VOLUME REQUIRED	/ IF SAMPLE COLLECTED	SAMPLE BOTTLE IDS
	SURFACE WATER	SEDIMENT				
☐ TCE VOC	☐	☒	☐	<u> </u>	☒	<u> </u>
☐ TCE SVCC	☐	☒	☐	<u> </u>	☒	<u> </u>
☐ TCE PESTICIDES	☐	☒	☐	<u> </u>	☒	<u> </u>
☐ TCE INORGANIC	☐	☒	☐	<u> </u>	☒	<u> </u>
☐ TCE METALS	☐	☒	☐	<u> </u>	☒	<u> </u>
☐ TCE/INORGANIC	☐	☒	☐	<u> </u>	☒	<u> </u>

NOTES/SKETCH

SAMPLE TAKEN AT EDGE OF ELEVATED VESSEL



SURFACE SOIL SAMPLE DATA RECORD

Project: ENRX, Inc. Site: Third Floor
 Project Number: 7170-40 Date: 10-5-94
 Sample Location ID: WT-106
 Time: Start: 0830 End: 0940 Signature of Sampler: P. Wilbur

SOIL SAMPLE

DEPTH OF SAMPLE 4-5" LOOSE
Pile on
Floor

EQUIPMENT USED FOR COLLECTION:

☐ HAND AUGER
☐ S.S. SPLIT SPOON
☐ SHOVEL
☒ HAND SPOON
☐ ALUMINUM PANS
☐ SS BUCKET

TYPE OF SAMPLE COLLECTED:

☒ DISCRETE
☐ COMPOSITE

SAMPLE OBSERVATIONS:

☐ ODOR
☐ COLOR BROWN, DRY

DECONTAMINATION FLUIDS USED:

☒ ALL USED
☐ ETHYL ALCOHOL
☐ 25% METHANOL/ 75% ASTM TYPE II WATER
☒ DEIONIZED WATER
☒ LIQUINOX SOLUTION
☐ HEXANE
☐ HNO₃ SOLUTION
☐ POTABLE WATER
☐ NONE

SOIL TYPE:

☐ CLAY
☐ SAND
☐ ORGANIC
☐ GRAVEL

SILTY, FINE DUST

FIELD GC DATA: ☐ FIELD DUPLICATE COLLECTED
 DUPLICATE ID _____

SAMPLE LOCATION SKETCH:

☒ YES see below
☐ NO

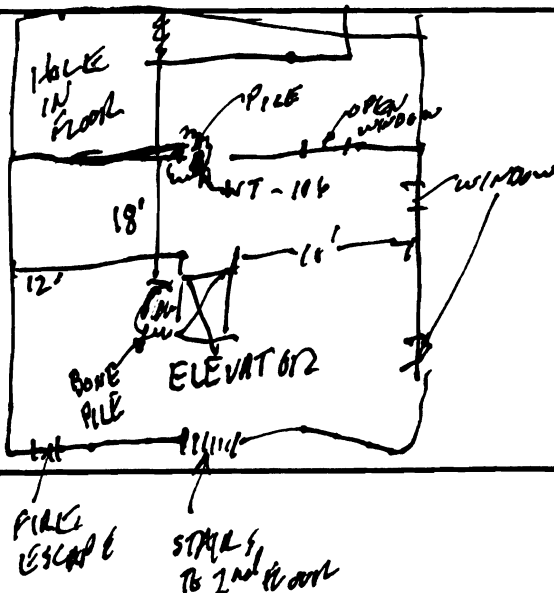
P10 = 0. ppm

O₂ = OPEN AIRWAYS

SAMPLES COLLECTED

✓ IF REQUIRED AT THIS LOCATION	MATRIX		✓ IF PRESERVED WITH ACID-BASE	VOLUME REQUIRED	✓ IF SAMPLE COLLECTED	SAMPLE BOTTLE IDS
	SURFACE WATER	SEDIMENT				
☒ TEL WCC	☐	☐	☐		☐	
☒ TEL SWCC	☐	☐	☐		☐	
☒ TEL Post/Rep	☐	☐	☐		☐	
☒ TEL incg	☐	☐	☐		☐	
☒ EPTX Methan	☐	☐	☐		☐	
☒ ignit/acc: fluid	☐	☐	☐		☐	

NOTES/SKETCH



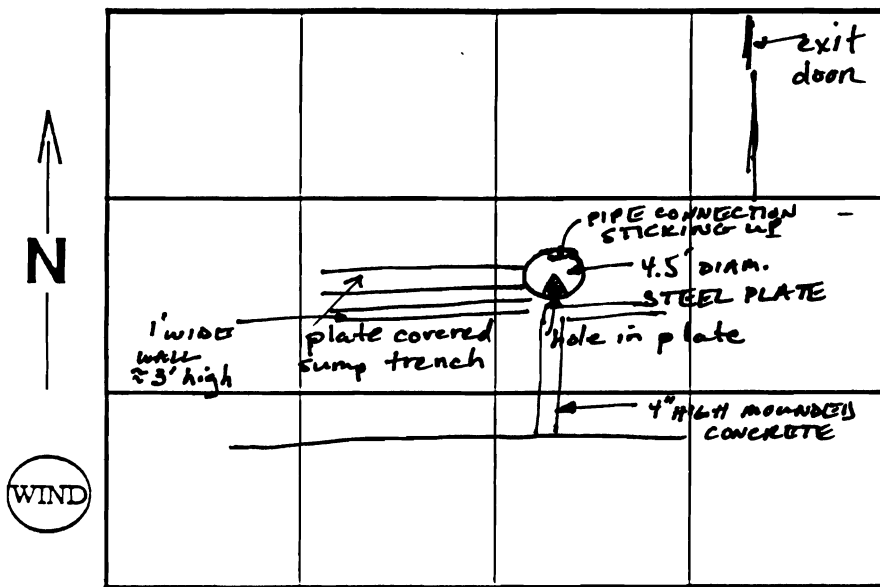
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SUMP/DRY WELL/STRUCTURE SAMPLING RECORD

PAGE 1 OF 3

SITE ENRY STRUCTURE TYPE O/W SEPARATOR
 STRUCTURE ID ENCD101 DATE 10-4-94 TIME 1030 END 1110
 COORDINATES _____

PLAN VIEW OF STRUCTURE SITE W/ DIMENSIONS



SCALE 1 INCH= NTS FEET

CREW MEMBERS:

1. D. LOVJOY
2. C. TALBOT
3. D. WILKIE
4. B. BUTLER
- 5.
- 6.

MONITORING EQUIPMENT:

PI Meter ☒ N
 Explosive Gas ☒ N
 Avail. Oxygen ☒ N
 OVA ☒ Y N
 Other B-K 1302 MULTIGAS
MONITOR

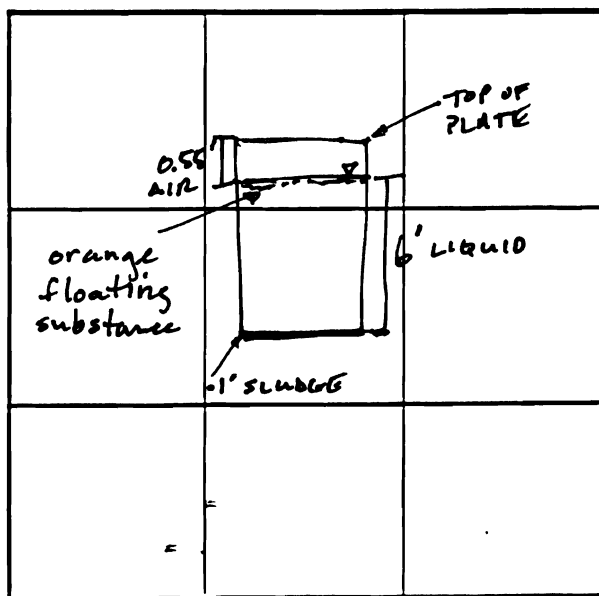
Photographs, Roll No

Exposure No

HEALTH AND SAFETY

Protection Level B
 Initial PI 0. ppm
 Initial LEL 1 %
 Initial O₂ 20.8 %

CROSS SECTION OF STRUCTURE



SCALE 1 INCH= NTS FEET

Logged by D Wilkie

Checked by B Butler

SUMP/DRY WELL/STRUCTURE SAMPLING RECORD

PAGE 2 OF 3

SITE ENRX STRUCTURE TYPE Oil Separator
STRUCTURE ID ENC101 DATE 10-4-94 TIME 1030 END 1110

LIQUID DATA:

Liquid Depth 6' ft. from top of tank Sample ID None Collected
Temperature 21.2 Degrees C. Sample Observations
pH 7.48 Units ☐ Odor None
Specific Conductivity 300 ~~umhos/cm~~ ms/cm ☐ Color Ben clear
☐ Layered orange floating scum
☐ (~.3-.4' thick)
PI Meter (Headspace) 25.3 ppm air between lid & water
Field GC Screening ☐ Yes ☒ No

Equipment Used for Collection None

Decontamination Fluids Used N/A

SLUDGE/SEDIMENT DATA:

Depth to Sediment 6.1 ft. from top of tank Sample ID ENC101XXX94XX
Depth to Structure Bottom _____ ft. from _____

Type of Sample Collected ☒ Discrete
☐ Composite

Sample Observations

☐ Odor _____
☐ Color gray-black
☐ _____
☐ _____

PI Meter (Headspace) 30 ppm
Field GC Screening ☐ Yes ☒ No

Equipment Used for Collection Bucket Auger

Decontamination Fluids Used Water/Liquinox, DI Water

ANALYTICAL PARAMETERS:

	LIQUID	SEDIMENT		LIQUID	SEDIMENT
☒ TCL VOC	☐	☒	☐ TCLP	☐	☐
☒ TCL SVOC	☐	☒	☒ TCL PEST/PCBs	☐	☒
☐ TPH	☐	☐	☒ EPTOX Metals	☐	☒
☒ TAL METALS	☐	☒	☒ Ign./React/Corr.	☐	☒

SUMP/DRY WELL/STRUCTURE SAMPLING RECORD

PAGE 3 OF 3

SITE ENRX STRUCTURE TYPE O/W Separator
STRUCTURE ID ENC101 DATE 10/4/94 TIME 1030 END 1110

COMMENTS: EVERYTIME SAMPLE WAS PULLED OUT, PID
READ 25-30 ppm.

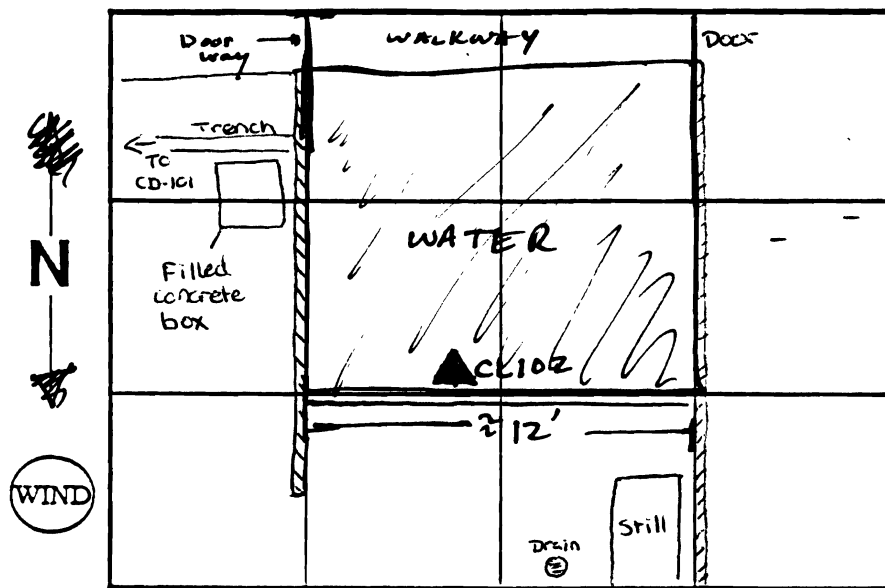
HEALTH AND SAFETY MONITORING NOTES: LEL/O₂ remained @ background.
All sampling performed @ Level B PPE.

SUMP/DRY WELL/STRUCTURE SAMPLING RECORD

PAGE 1 OF 3

SITE ENRY STRUCTURE TYPE Collection Pool
 STRUCTURE ID ENCL102 DATE 10/4/99 TIME 1145 END 1220
 COORDINATES _____

PLAN VIEW OF STRUCTURE SITE W/ DIMENSIONS



SCALE 1 INCH = NTS FEET

CREW MEMBERS:

1. P. Lovejoy
2. C. TALBOT
3. D. WILKIE
4. B. BUTLER
5. _____
6. _____

MONITORING EQUIPMENT:

PI Meter Y N
 Explosive Gas Y N
 Avail. Oxygen Y N
 OVA Y N
 Other Brüel and Kjaer
1302 Multigas Monitor

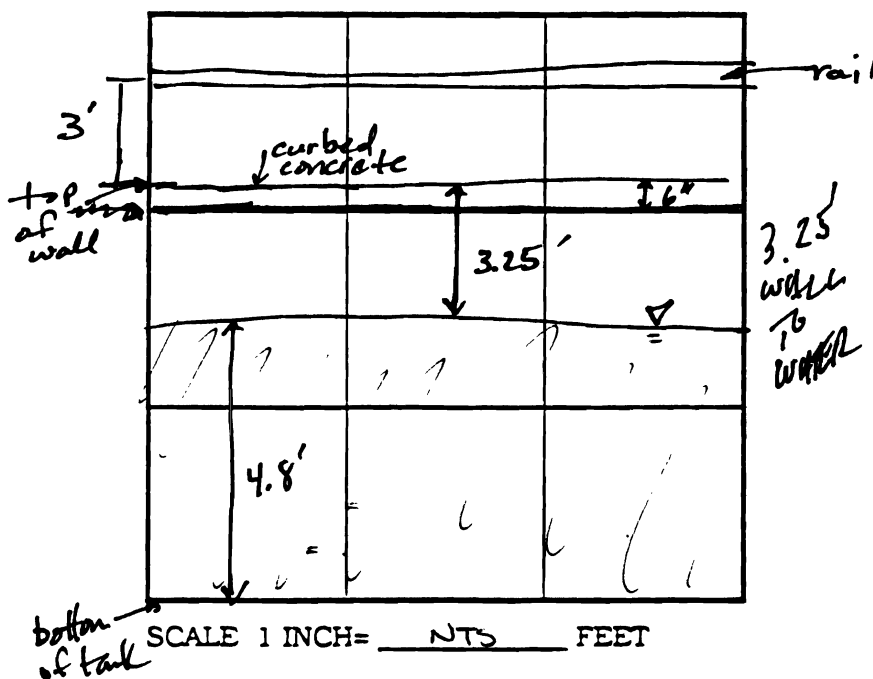
Photographs, Roll N/A

Exposure N/A

HEALTH AND SAFETY

Protection Level B
 Initial PI 0 ppm
 Initial LEL 1 %
 Initial O₂ 20.6 %

CROSS SECTION OF STRUCTURE



SCALE 1 INCH = NTS FEET

Logged by D Wilkie

Checked by B Butler

SUMP/DRY WELL/STRUCTURE SAMPLING RECORD

PAGE 2 OF 3

SITE ENRX STRUCTURE TYPE Collection Pool
 STRUCTURE ID ENCL102 DATE 10-4-94 TIME 1145 END 1220

LIQUID DATA:

Liquid Depth 4.8' ft. from N. edge concrete Sample ID ENCL102XXX94XX
 Temperature 10.2 Degrees C. Sample Observations
 pH 6.92 Units ☐ Odor NO
 Specific Conductivity 461 ~~umhos/cm~~ MS/cm ☐ Color clear, some turbidity
☐ Layered NO
☐
☐

PI Meter (Headspace) 0 ppm

Field GC Screening ☐ Yes ☒ No

Equipment Used for Collection Pak-bomb sampler /w teflon cord.

Decontamination Fluids Used Liquinox / Di Water

SLUDGE/SEDIMENT DATA:

Depth to Sediment N/A ft. from N/A Sample ID N/A
 Depth to Structure Bottom _____ ft. from _____

Type of Sample Collected ☐ Discrete
☐ Composite

Sample Observations

☐ Odor _____
☐ Color _____
☐ _____
☐ _____

PI Meter (Headspace) _____ ppm

Field GC Screening ☐ Yes ☐ No

Equipment Used for Collection N/A

Decontamination Fluids Used N/A

ANALYTICAL PARAMETERS:

	LIQUID	SEDIMENT		LIQUID	SEDIMENT
☒ TCL VOC	☒	☐	☐ TCLP	☐	☐
☒ TCL SVOC	☒	☐	☒ TCL PEST/PCBs	☒	☐
☐ TPH	☐	☐	☒ EPTOX Metals	☒	☐
☒ TAL METALS	☒	☐	☒ Ign./React/Corr.	☒	☐

SUMP/DRY WELL/STRUCTURE SAMPLING RECORD

PAGE 3 OF 3

SITE ENRX STRUCTURE TYPE Collection Pool
STRUCTURE ID ENCL102 DATE 10-4-94 TIME 1145 END 1220

COMMENTS: Sampling performed @ level B

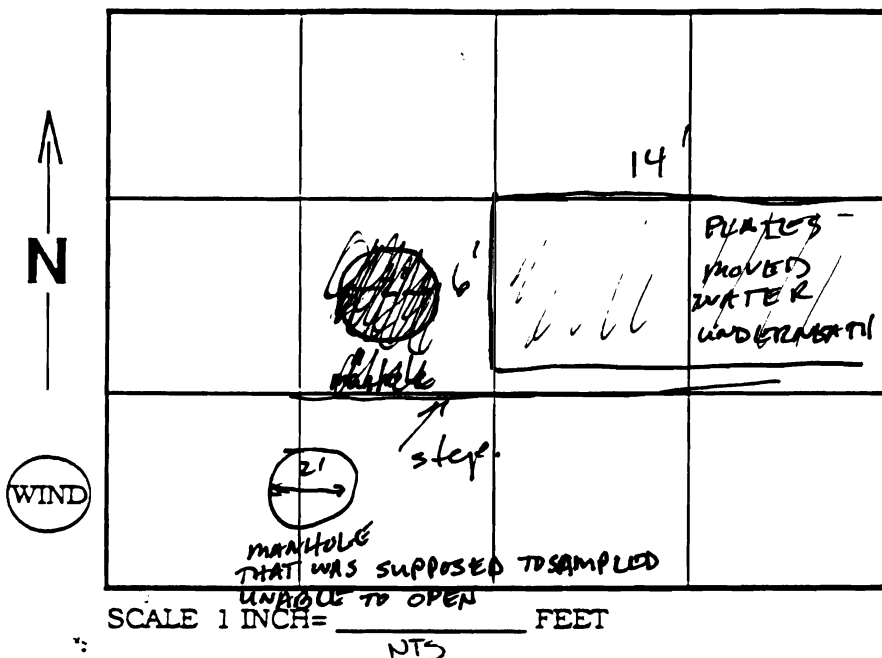
HEALTH AND SAFETY MONITORING NOTES: LEL/O₂ remained @ background

SUMP/DRY WELL/STRUCTURE SAMPLING RECORD

PAGE 1 OF 3

SITE ENRY STRUCTURE TYPE OW Separator
 STRUCTURE ID ENCD103/CH03 DATE 10-4-99 TIME 1230 END 1550
 COORDINATES _____

PLAN VIEW OF STRUCTURE SITE W/ DIMENSIONS



CREW MEMBERS:

1. D. Lovejoy
2. C. TALBOT
3. D. WILKIE
4. B. BUTLER
- 5.
- 6.

MONITORING EQUIPMENT:

PI Meter ☒ N
 Explosive Gas ☒ N
 Avail. Oxygen ☒ N
 OVA ☒ N
 Other Briel and Kjaer
1302 Multigas Monitor

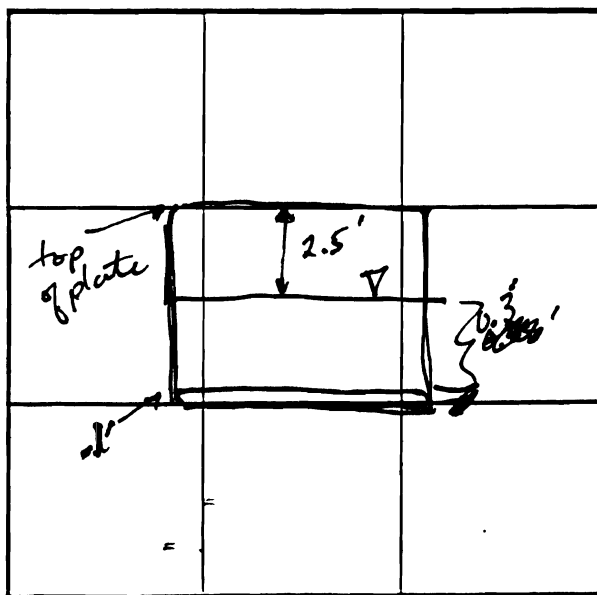
Photographs, Roll N/A

Exposure N/A

HEALTH AND SAFETY

Protection Level B
 Initial PI 0 ppm
 Initial LEL — %
 Initial O₂ 20.6 %

CROSS SECTION OF STRUCTURE



SCALE 1 INCH= NTS FEET

Logged by D. Wilkie

Checked by B Butler

SUMP/DRY WELL/STRUCTURE SAMPLING RECORD

PAGE 2 OF 3

SITE ENRX STRUCTURE TYPE o/w separator
STRUCTURE ID ENCL103/CU03 DATE 10/4/94 TIME 1230 END 1530

LIQUID DATA:

Liquid Depth ^{BB 10/5/94} 0.5 2.5 ft. from Floor Sample ID ENCL103XXX94XX
Temperature N/A Degrees C. Sample Observations
pH N/A Units ☐ Odor
Specific Conductivity N/A μ mhos/cm ☐ Color
☐ Layered
☒ clear to rusty
☐

PI Meter (Headspace) 0 ppm
Field GC Screening ☐ Yes ☒ No

Equipment Used for Collection ^{BB 10/5/94} Liquinox, DI water Filled Directly

Decontamination Fluids Used Liquinox, DI water

SLUDGE/SEDIMENT DATA:

Depth to Sediment ^{BB 10/5/94} 2.0 2.0 ft. from Floor Sample ID ENCL103XXX94XX
Depth to Structure Bottom ^{BB 10/5/94} 3.0 3.0 ft. from Floor

Type of Sample Collected ☒ Discrete
☐ Composite

Sample Observations
☐ Odor
☒ Color Brown w/ orange Flcks.
☐
☐

PI Meter (Headspace) 0 ppm
Field GC Screening ☐ Yes ☒ No

Equipment Used for Collection Bucket Auger

Decontamination Fluids Used Liquinox, DI water

ANALYTICAL PARAMETERS:

	LIQUID	SEDIMENT		LIQUID	SEDIMENT
☒ TCL VOC	☒	☒	☐ TCLP	☐	☐
☒ TCL SVOC	☒	☒	☒ TCL PEST/PCBs	☒	☒
☐ TPH	☐	☐	☒ EPTOX Metals	☒	☒
☒ TAL METALS	☒	☒	☒ Ign./react/corr.	☒	☒

SUMP/DRY WELL/STRUCTURE SAMPLING RECORD

PAGE 3 OF 3

SITE ENRX STRUCTURE TYPE O/W Separator
STRUCTURE ID ENC0103/CL103 DATE 10/4/94 TIME 1230 END 1550

COMMENTS: Sample collected from Pit instead of sewer Manhole cover. Could not open manhole for access to original location planned.

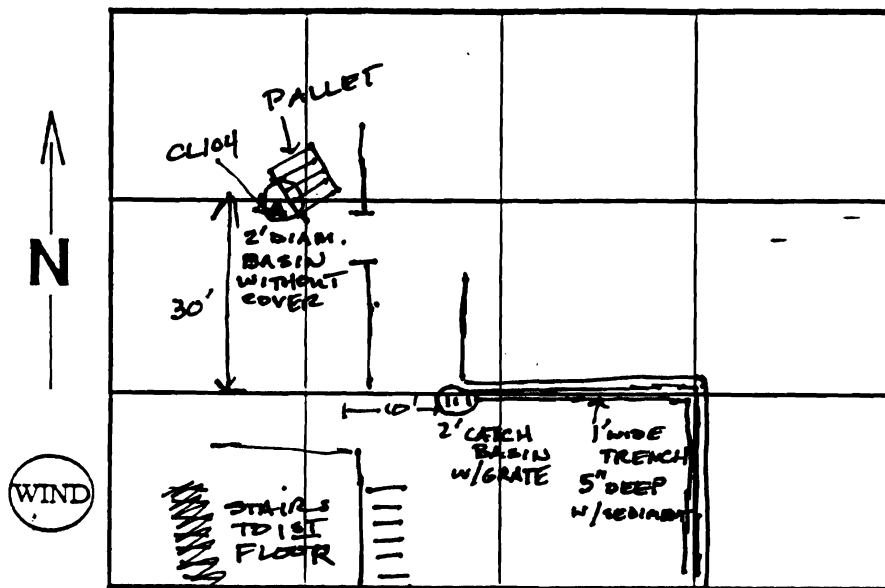
HEALTH AND SAFETY MONITORING NOTES: Sample collected @
Level B PPE. See B+K 1302 data. Max. VOC compound
concentrations in air @ truck bay where samples
collected:
1,2-DCE = 2.81 ppm
1,1,1-TCA = 0.349 ppm ^{0.510 ppm} 1.0 ppm
1,1-DCE = 0.642 ppm
CCl₄ = 0.151 ppm

SUMP/DRY WELL/STRUCTURE SAMPLING RECORD

PAGE 1 OF 1

SITE ENRX, INC. STRUCTURE TYPE Catch basin
 STRUCTURE ID CL104 DATE 10-5-94 TIME 0955 END 1010
 COORDINATES _____

PLAN VIEW OF STRUCTURE SITE W/ DIMENSIONS



SCALE 1 INCH = NTS FEET

CREW MEMBERS:

1. D. Lovejoy
2. C. TALBOT
3. D. WILKIE
4. B. BUTLER
- 5.
- 6.

MONITORING EQUIPMENT:

PI Meter (Y) N
 Explosive Gas (Y) N
 Avail. Oxygen (Y) N
 OVA Y N
 Other _____

Photographs, Roll _____

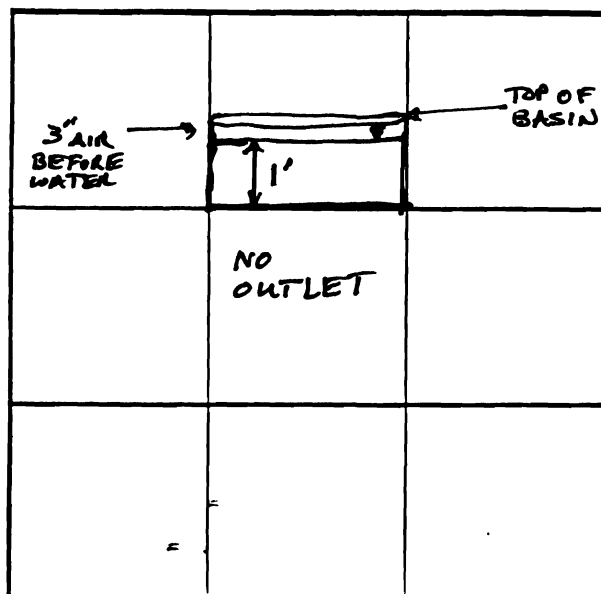
Exposure _____

HEALTH AND SAFETY

Protection Level B
 Initial PI 765 ppm
 Initial LEL 1 %
 Initial O₂ 20.3 %

WATER: COLOR: GRAY

CROSS SECTION OF STRUCTURE



SCALE 1 INCH = NTS FEET

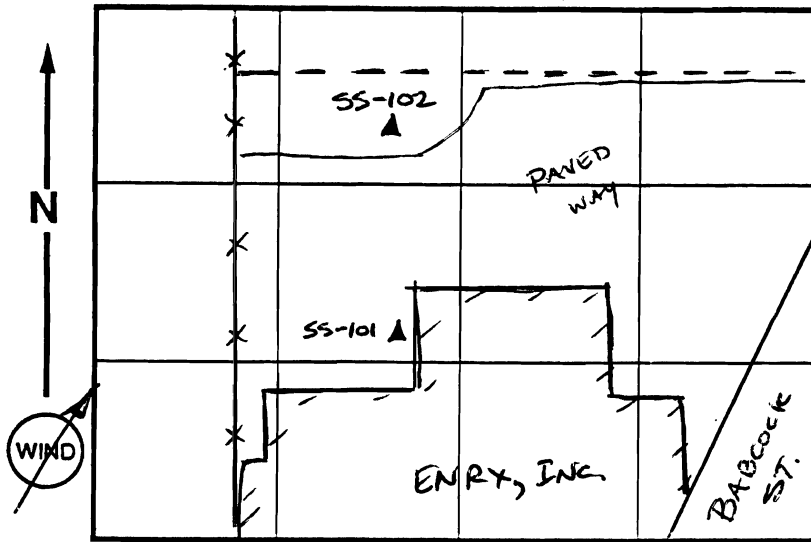
Logged by D. Wilkie

Checked by B. Butler

SURFACE SOIL SAMPLING RECORD

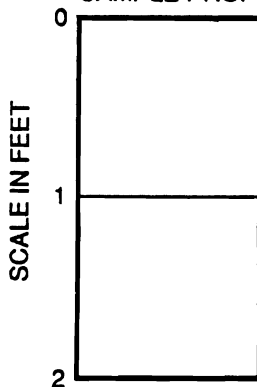
Site: ENRY, INC. Project No. 7170-40
 Location No. SS-102 Date 10-28-94 Time 10:15 End 10:20
 Coordinates _____ AOC ENRY, INC.

SKETCH MAP OF SAMPLING SITE



SCALE 1" = ~ 50 FT.

SAMPLE PROFILE



No.	Sample No.	Depth (ft.)
S-1	ENSS102XXXXXX	0.5-1.0
S-2	ENSS102XXXXXX	0.5-1.0
S-3		

Sampling Equipment:

SS. Spoon & Bucket

Decon. Materials:

LIQUINOX & Pot. Water
DI Water

Crew Members:

1. Tom Lowkey
2. Ashley Foster
- 3.
- 4.
- 5.
- 6.

Monitor Equipment:

PI Meter (Y) N
 Explosive Gas Y N
 Avail. Oxygen Y N
 OVA Y N
 Other _____

Photographs: (Roll Exposure)

SAMPLE DESCRIPTION:

Black, ORGANIC, FILL - GRAVELLY
SAND & SILT
FROM GRASSY, UNPAVED AREA

NOTES: Collected Sample, Dup, MS/MSD

3 4-oz. VOA JARS TEL VOA
6 8-oz. SVOA JARS TEL SVOA
" P/PCB
" Inorganics
EP TOX

References:

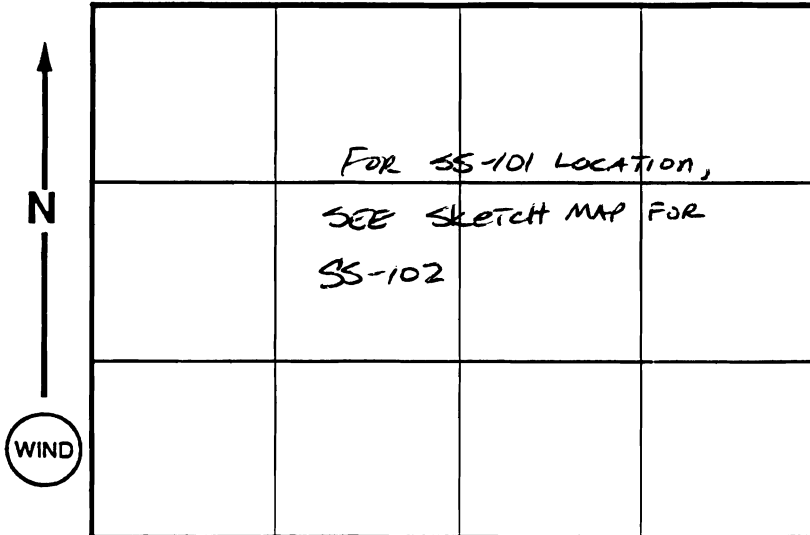
Field Book #: FIELD Sampling
 Page #: 30
 Attachments:

Signature: John D. Lowkey

SURFACE SOIL SAMPLING RECORD

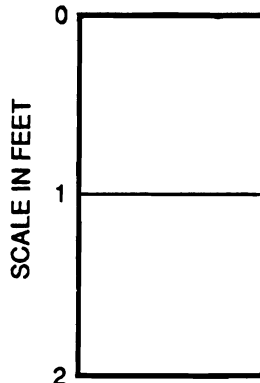
Site: ENRX, INC. Project No. 7170-40
 Location No. SS-101 Date 10-28-94 Time 10:30 End 10:35
 Coordinates _____ AOC ENRX, INC.

SKETCH MAP OF SAMPLING SITE



SCALE 1" = _____ FT.

SAMPLE PROFILE



No.	Sample No.	Depth (ft.)
S-1	ENSS101XX194X	0.5-1.0
S-2		
S-3		

Sampling Equipment:

S.S. Spoon & Bucket

Decon. Materials:

LIQUINOX & Pot. Water,
DI water

Crew Members:

1. Tom Longley
2. ASHLEY FOSTER
3. _____
4. _____
5. _____
6. _____

Monitor Equipment:

PI Meter Y N
 Explosive Gas Y N
 Avail. Oxygen Y N
 OVA Y N
 Other _____

Photographs: (Roll Exposure)

SAMPLE DESCRIPTION:

Black, Dry, FILL

NOTES:

ADJACENT TO ENRX BLDG. ON
NW-SIDE - MAY contain
Bitumen pieces FROM TAR
DRIVEWAY 1- 4oz. VOA TAR
2- 8oz. SVOA TARS

References:

Field Book #: FIELD SAMPLING
 Page #: 30
 Attachments: _____

Signature:

John D. Longley

SECTION 2.0

BORING LOGS AND MONITORING WELL INSTALLATION DIAGRAMS

2018

2019

2020

2021

2022

2023

2024

2025

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Test Boring Log

Project ENRX, INC.		Boring/Well No. MW-101	Project No. 7170-40	
Client NYSDEC	Site ENRX, INC.		Sheet No. <u>1</u> of <u>1</u>	
Logged By T. LONGLEY		Ground Elevation	Start Date 10-27-94	Finish Date 10-27-94
Drilling Contractor A.D.I.		Driller's Name MARK SEILER	Rig Type MOBILE ATV	
Drilling Method H.S.A./AIR ROTARY		Protection Level C DERMAL	P.I.D. (eV) 10.0	Casing Size 4.25" Auger Size I.D.
Soil Drilled 7.5'	Rock Drilled 7.5'	Total Depth 15'	Depth to Groundwater/Date 6.23' BELOW TOP PVC 10/28/94	
		Piez ☐ Well ☒ Boring ☐		

Depth(Feet)	Sample No. & Penetration/ Recovery (Feet)	Sample Type	SPT Blows/6" or Core Rec./Rqd. %	SPT-N (Blows/Ft.)	Graphic Log	Sample Description	USCS Group Symbol	Notes on Drilling	Monitoring		Lab Tests
									(ppm)		
									PI Meter Field Scan	PI Meter Head Space	
1	S-1 2.0 1.4	SPT	12- 13- 19- 12	32		GRAVELLY SAND-FILL-WET AT SURFACE, BRICK PIECES, TAR PIECES OVER TAN GRAVELLY FINE SAND	Fill	0-4' Rough THEN EASY FROM 4' TO 7.5'	* 4.5	NR	
3	S-2 2.0 0.6	SPT	3- 3- 3- 3	6		BLACK, GRAVELLY, SANDY SILT TO SILTY SAND, BRICK PIECES, MOIST	Fill		Bkg.	NR	
5	S-3 2.0 0.3	SPT	3- 2- 7- 3	9		BROWN, WET FILL, GRAVELLY SANDY SILT - WOOD PIECE (R.R. TIE?)	Fill		Bkg.	NR	
7	S-4 1.5 0.3	SPT	3- 6- 50/4"	>56		AS ABOVE THEN LAST 0.2' SATURATED BLACK SILTY SAND SANDY SILT W/ ROCK & WOOD BEDROCK AT 7.5'	Fill		Bkg.	NR	ENB3101 XX694XX
9						DRILLED W/ AIR ROTARY USING 3 7/8" ROLLER BIT FROM 7.5 TO 15' USING AUGERS AS CASING. DRILLED FROM 17:15 → 18:40 W/ AIR ROTARY APPROX. 12 MIN./FOOT	Rock				
11											
13											
15						B.O.B. 15'					

ABB Environmental Services

ABB Environmental Services

Test Boring Log

Project ENRX, INC.				Boring/Well No. MW-102		Project No. 7170-40	
Client NYSDEC			Site ENRX, INC.			Sheet No. <u>1</u> of <u>1</u>	
Logged By T. LONGLEY			Ground Elevation		Start Date 10-28-94		Finish Date 10-28-94
Drilling Contractor A.D.I.			Driller's Name MARK SEILER			Rig Type MOBILE ATV.	
Drilling Method H.S.A. / AIR ROTARY			Protection Level C DERM.		P.I.D. (eV) 10.0	Casing Size 4.25"	Auger Size I.D.
Soil Drilled 8'		Rock Drilled 7'		Total Depth 15'	Depth to Groundwater/Date 6.1' BELOW TOP PVC 10/31/94		Piez ☐ Well ☒ Boring ☐

Depth (Feet)	Sample No. & Penetration/ Recovery (Feet)	Sample Type	SPT Blows/6" or Core Rec./Rqd. %	SPT-N (Blows/Ft.)	Graphic Log	Sample Description	USCS Group Symbol	Notes on Drilling	Monitoring (ppm)		Lab Tests
									PI Meter Field Scan	PI Meter Head Space	
1	S-1 2.0 1.8	SPT	5-6-7-5	13		BLACK, SOOT-LIKE FILL, GRAVELLY SANDY SILT, BRICKS, COAL, ASH, SLAG, MED. DENSE, DRY TO DAMP	FILL		0	NR	
3	S-2 2.0 1.3	SPT	10-9-14-36	23		AS ABOVE BUT MORE VARIABLE IN COLOR - CINDERS, ASH, BRICK, COAL, MED. DENSE, DRY TO DAMP	FILL		0	NR	
5	S-3 2.0 1.2	SPT	8-5-5-8	10		FILL AS ABOVE TO 0.5' 0.5'-1.2' GRAY SILT W/ GRAVEL & DEBRIS, STIFF, DRY TO DAMP	FILL		0	NR	
7	S-4 2.0 1.2	SPT	9-15-30-50 1/4"	45		0-0.7' FILL AS ABOVE 0.7-1.2' BROWN SILT & SAND, WET, EVENLY GRADED	FILL SM		0	NR	ENBS102 XX694XX
9						BEDROCK AT 8'					
11						DRILLED W/ AIR ROTARY FROM 9:35 TO 10:55 APPROX. 11 MIN./FOOT - DRY ROCK - MUCH DUST, POSSIBLY FEW FRACTURES					
13											
15						B.O.B. 15'					

ABB Environmental Services

Test Boring Log

Project ENRX, INC.		Boring/Well No. MW-103	Project No. 7170-40	
Client NYSDEC	Site ENRX, INC.		Sheet No. <u>1</u> of <u>1</u>	
Logged By T. LONGLEY	Ground Elevation	Start Date 10-28-94	Finish Date 10-28-94	
Drilling Contractor A.D.I.	Driller's Name MARK SEILER		Rig Type MOBILE ATV	
Drilling Method H.S.A.	Protection Level C DERMAL	P.I.D. (eV) 10.0	Casing Size N/A	Auger Size 4.25" I.D.
Soil Drilled 8'	Rock Drilled —	Total Depth 8'	Depth to Groundwater/Date 2.64' BELOW PVC	
		Piez ☐ Well ☒ Boring ☐		

Depth(Feet)	Sample No. & Penetration/ Recovery (Feet)	Sample Type	SPT Blows/6" or Core Rec./Rqd. %	SPT-N (Blows/Ft.)	Graphic Log	Sample Description	USCS Group Symbol	Notes on Drilling	Monitoring		Lab Tests
									(ppm)		
									PI Meter Field Scan	PI Meter Head Space	
1	S-1 2.0 0.8	SPT	30- 28- 30- 13	58		BLACK GRAVELLY SILT, SAND, COARSE CINDERS, COAL - DRY - BRICKS	FILL		Bkg.	NR	
3	S-2 3.0 0.2	SPT	7- 5- 3- 8	8		STINKS LIKE SEWER - BRICKS - RUBBER TIRES SATURATED * 9ppm OVER H.S.A.	FILL		Bkg. *	NR	
5	S-3 2.0 0.1	SPT	8- 7- 5- 2	12		BRICK PIECES - V. LITTLE RECOVERY HIT REFUSAL AT 6' - MOVE OVER & DRILL TO 8'	FILL		Bkg.	NR	
7	S-4 2.0 0.0	SPT	1- 1- 1- 1	2		NO RECOVERY - DRILLER SAYS THIS IS BEDROCK WHERE 6' DEPTH WAS NOT	FILL		NR	NR	
						B.O.B. @ 8' BGS NO ANALYTICAL SAMPLES COLLECTED AT THIS BORING LOCATION					

ABB Environmental Services

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BEDROCK MONITORING WELL CONSTRUCTION DIAGRAM

Project ENRY, INC. Study Area _____ Driller A.D.I.
 Project No. 7170-40 Boring No. MW-101 Drilling Method 4.25" HSA ± 3 3/8" AIR ROTARY
 Date Installed 10-27-94 Development Method WATER
 Field Geologist Tan Longley

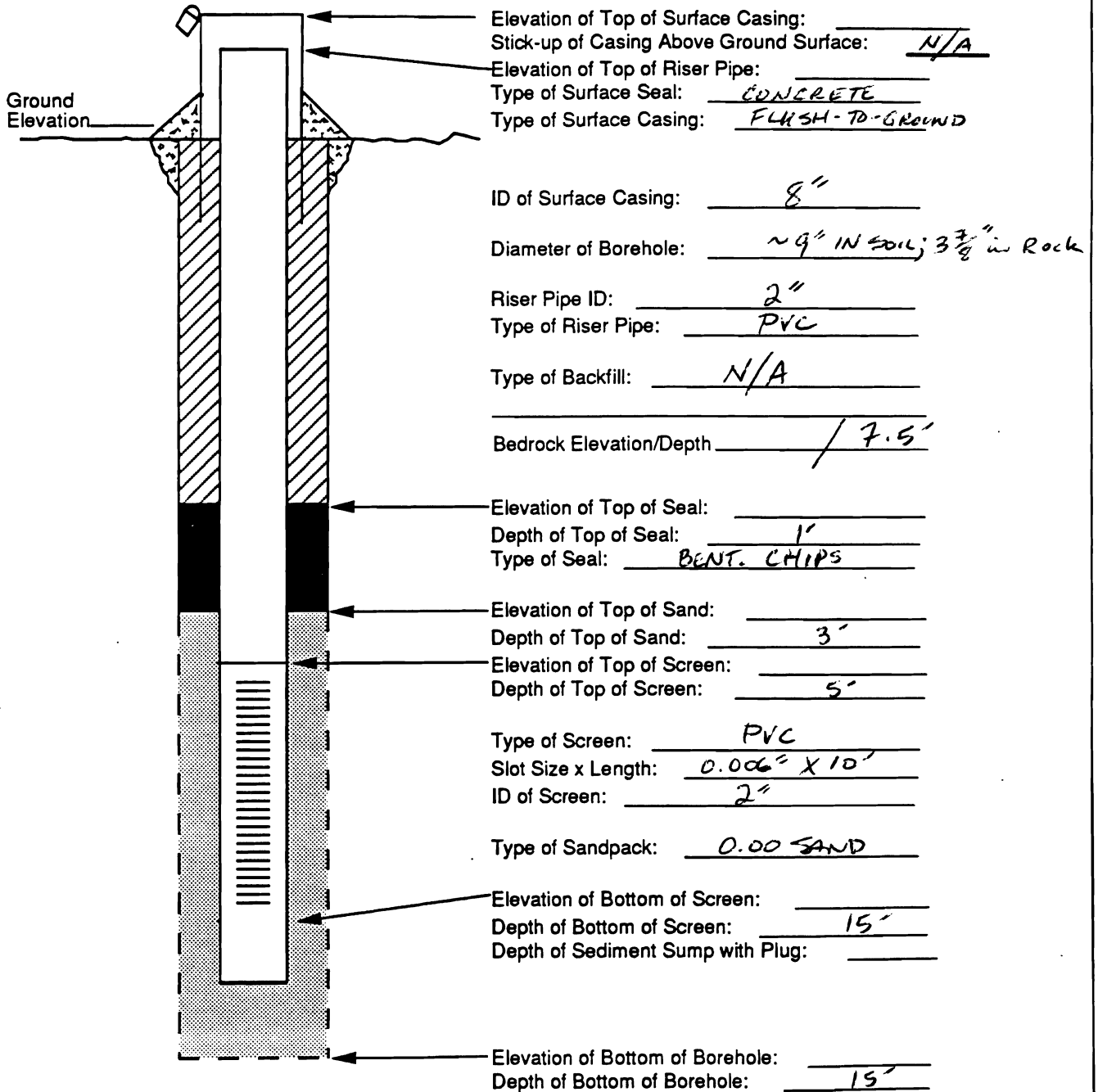


FIGURE 4-12
BEDROCK MONITORING WELL CONSTRUCTION DIAGRAM
NYSDEC QUALITY ASSURANCE PROGRAM PLAN

ABB Environmental Services, Inc.

BEDROCK MONITORING WELL CONSTRUCTION DIAGRAM

Project ENRY, INC. Study Area _____ Driller A.D.I.
 Project No. 7170-40 Boring No. MW-102 Drilling Method 4.25" H.S.A. / 3 3/8" AIR ROTARY
 Date Installed 10-27-94 Development Method WATERAA
 Field Geologist Tom Longley

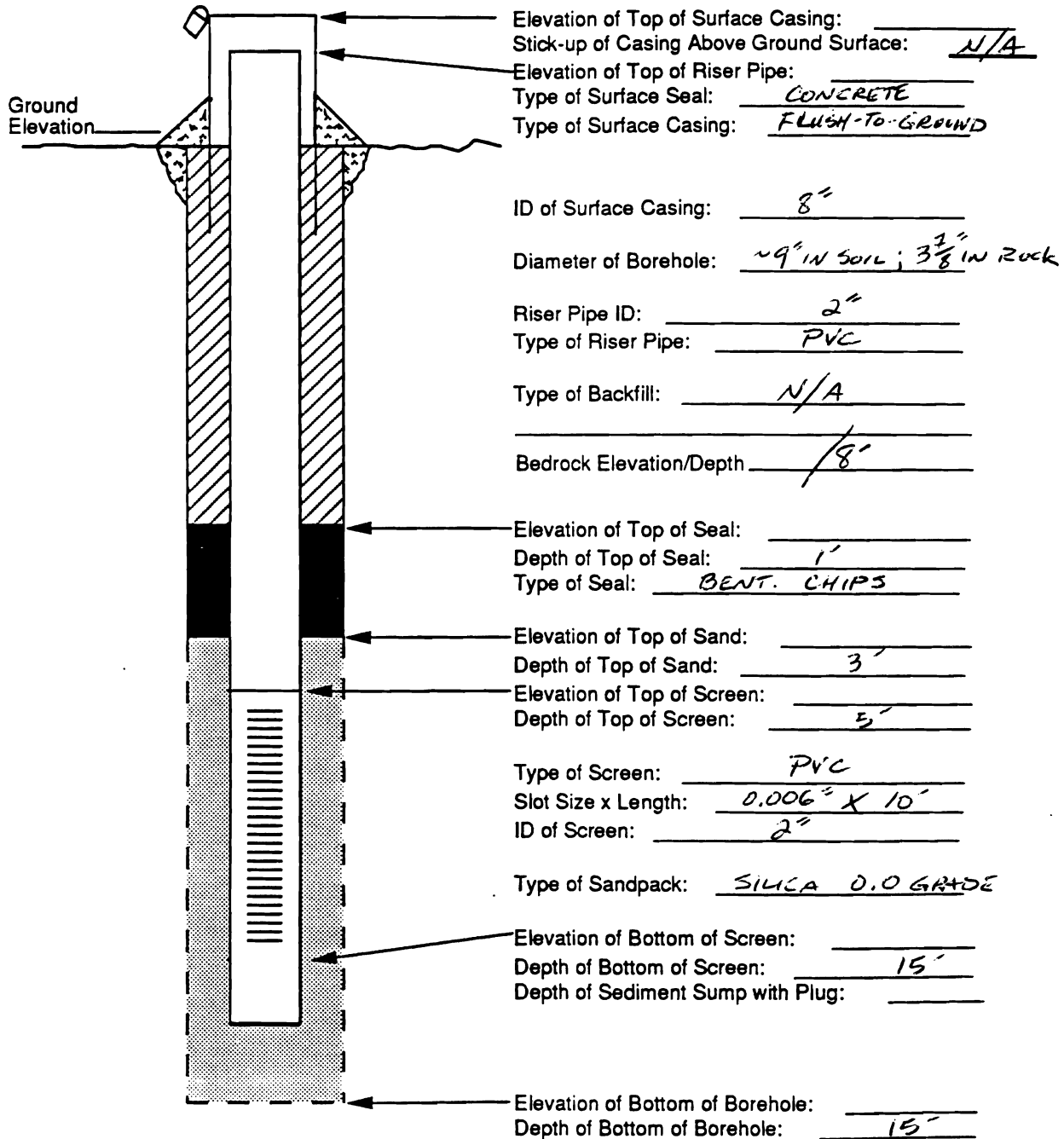


FIGURE 4-12
BEDROCK MONITORING WELL CONSTRUCTION DIAGRAM
NYSDEC QUALITY ASSURANCE PROGRAM PLAN

ABB Environmental Services, Inc.

OVERBURDEN MONITORING WELL CONSTRUCTION DIAGRAM

Project ENRY, INC. Study Area _____ Driller A.D.I.
 Project No. 7170-40 Boring No. MW-103 Drilling Method 4.25" H.S.A.
 Date Installed 10-28-94 Development Method CENT. PUMP
 Field Geologist Tom Longley

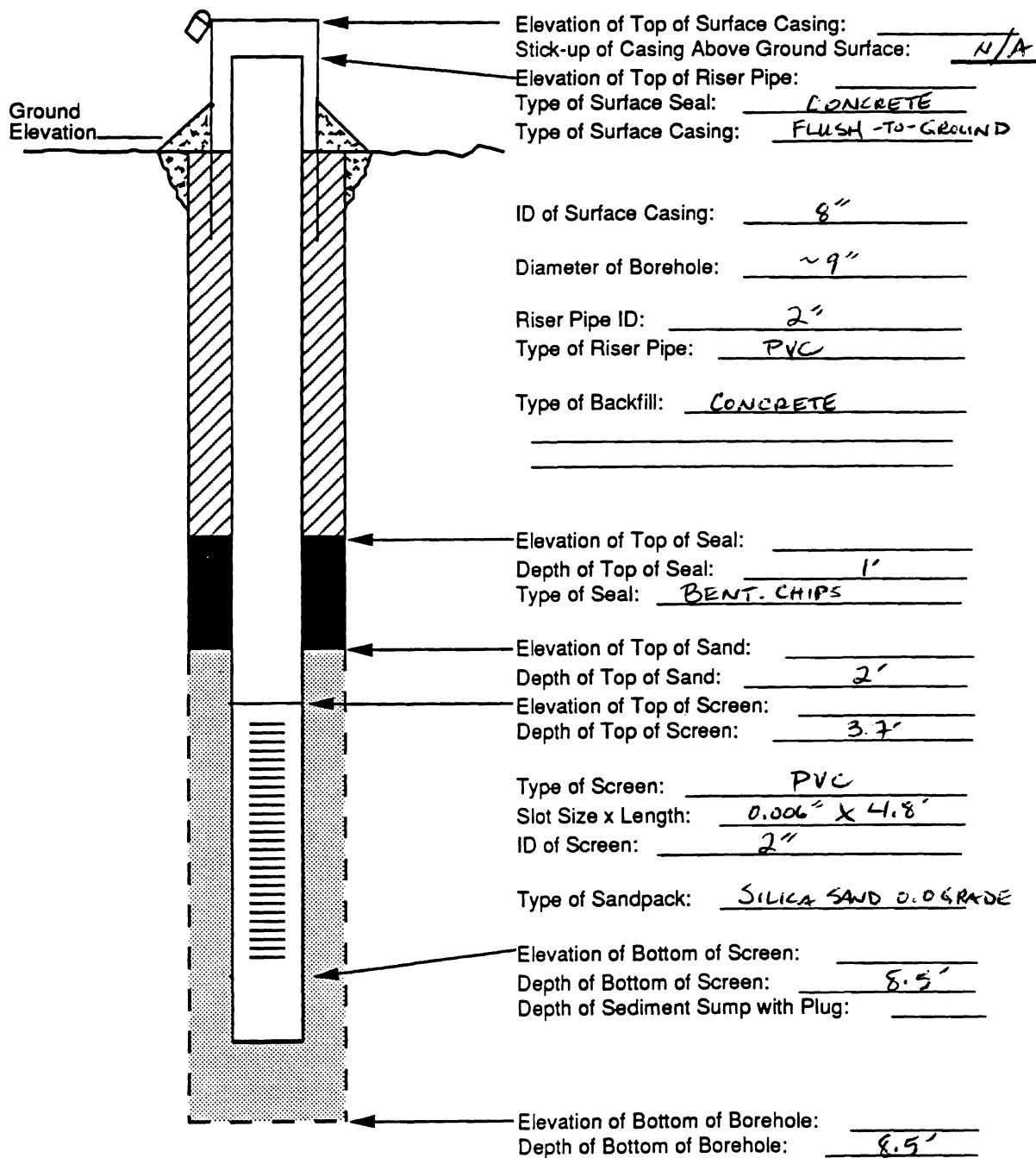


FIGURE 4-11
OVERBURDEN MONITORING WELL CONSTRUCTION DIAGRAM
NYSDEC QUALITY ASSURANCE PROGRAM PLAN

ABB Environmental Services, Inc.

WELL DEVELOPMENT RECORD

Project: ENRY, INC.	Well Installation Date: 10-27-94	Project No. 7170-40
Client: NYSDEC	Well Development Date: 10-31-94	Logged by: TDL
Well/Site I.D.: MW-101	Weather: RAIN, CALM, 60's	Start Date: 10-31-94
Initial Water Level (ft): 6.4' Below TOP PVC	Start Time: 12:22	Finish Date: 10-31-94
Water Level during Initial Pumping/Purging (ft): N/A		Finish Time: 14:30
Water Level at Termination of Pumping/Purging (ft): N/A		

TOTAL	^{GALLONS} Number of Well Volumes	TIME	TEMP.	pH	Conductivity	Approximate Pumping Rate (gal/min)	Turbidity (NTU's)
	3.5	1240					
	5	1250	15.1				
	5.5	1255	15.1/15.2	6.94	1.61	0.07	7999
	7	1305	15.2/15.0	7.00	1.60	0.08	7999
	10	1325	15.0/15.1	6.9	1.61	0.07	897
	15	1430	15.1	7.0	1.61	0.07	570
	16.5	1430	14.8	7.0	1.64	0.07	520

% Solids

NOTES:

Developed using Watera Pump & hand-pumped to develop this well.

Considered developed based on stabilization of parameters.

Well Developer's Signature

FIGURE 4-13
WELL DEVELOPMENT RECORD
NYSDEC QUALITY ASSURANCE PROGRAM PLAN

ABB Environmental Services, Inc.

WELL DEVELOPMENT RECORD

Project: ENRY, INC.	Well Installation Date: 10-28-94	Project No. 7170-40
Client: NYSDEC	Well Development Date: 10-31-94 / 11-1-94	Logged by: TDL
Well/Site I.D.: MM-102	Weather: RAIN, CALM, 60's	Start Date: 10-31-94
Initial Water Level (ft): 6.1' Below TOP PVC		Finish Date: 11-1-94
Water Level during Initial Pumping/Purging (ft): N/A		Start Time: 15:35 10/31
Water Level at Termination of Pumping/Purging (ft): N/A		Finish Time: 08:10 11/1

TOTAL Number of Well Volumes	TIME	TEMP.	pH	Conductivity	Approximate Pumping Rate (gal/min) % Salinity	Turbidity (NTU's)
1.5	15:32					
Dry @ 3.0	10/31 16:30	12.2	7.08	2.41	0.11	163
4	11/1 08:00	12.4	7.1	2.22	0.10	445
Dry @ 5	08:10	12.1	7.14	2.27	0.10	480

NOTES:

Consider this well as developed. Used Wattera 3 hand bailed this well.

Well Developer's Signature

John D. Zyglis

FIGURE 4-13
WELL DEVELOPMENT RECORD
NYSDEC QUALITY ASSURANCE PROGRAM PLAN

ABB Environmental Services, Inc.

WELL DEVELOPMENT RECORD

Project: ENRY, INC.	Well Installation Date: 10-28-94	Project No. 7170-40	
Client: NYSDEC	Well Development Date: 10-31-94	Logged by: TDL	Checked by:
Well/Site I.D.: MW-103	Weather: RAIN, CALM, 60's	Start Date: 10-31-94	Finish Date: 10-31-94
Initial Water Level (ft): 2.64' Below TOP PVC	Start Time: 14:45	Finish Time: 15:15	
Water Level during Initial Pumping/Purging (ft): N/A			
Water Level at Termination of Pumping/Purging (ft): N/A			

TOTAL ^{GALLONS} Number of Well Volumes	TIME	TEMP.	pH	Conductivity	Approximate Pumping Rate (gpm/min) % SALINITY	Turbidity (NTU's)
10	14:55					
15	14:58	13.6	7.39	0.98	0.04	>999
20	15:02					
22	15:05	13.5	7.13	1.04	0.04	>999
25	15:06					
30	15:10	13.7	7.19	1.05	0.04	>999
35	15:15	13.3	7.15	1.05	0.04	799

NOTES:

STILL VERY DIRTY AFTER 35 GALS. - Slight sheen on devel. water.
 Used centrifugal pump w/ Foot valve/hose assembly inserted into well for development purposes.

Based on stabilization, consider this well developed.

Well Developer's Signature

John D. Zyl

FIGURE 4-13
 WELL DEVELOPMENT RECORD
 NYSDEC QUALITY ASSURANCE PROGRAM PLAN

ABB Environmental Services, Inc.

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GROUNDWATER SAMPLE FIELD DATA RECORD

Project: ENRX, INC
Project Number: 7170-40

Site: MW-101
Date: 11/30/94

Sample Location ID: ENMW101XXXX94XX

Time: Start: 0715 End: 0750

Signature of Sampler: R-K Buttl

Water Level/Well Data

Well Depth 14.5 Ft. ☒ Measured ☐ Historical ☒ Top of Well ☐ Top of Protective Casing

Well Riser Stick-up -0.60 Ft. (from ground) (Flush Mt) Protective - 0.6 Ft. Casing/Well Difference

Protective - 0 Ft. Casing

Depth to Water 5.96 Ft. (5.91 on 11/29/94) Well Material: ☒ PVC ☐ SS Well Locked?: ☒ Yes ☐ No Well Dia. ☒ 2 inch ☐ 4 inch ☐ 6 inch

Water Level Equip. Used: ☒ Elect. Cond. Probe ☐ Float Activated ☐ Press. Transducer

Height of Water Column 8.54 Ft. ☒ .16 Gal/Ft. (2 in.) ☐ .65 Gal/Ft. (4 in.) ☐ 1.5 Gal/Ft. (6 in.) ☐ Gal/Ft. (in.)

1.4 Gal/Vol. 4 Total Gal Purged

Well Integrity: Prot. Casing Secure ☒ Yes ☐ No Concrete Collar Intact ☒ Yes ☐ No Other ☐ Yes ☐ No

Equipment Documentation

Purging/Sampling Equipment Used:

Decontamination Fluids Used:

(✓ If Used For)

Purging	Sampling	Equipment ID
☒	☒	Peristaltic Pump
☐	☒	Submersible Pump
☐	☒	Bailer
☐	☒	PVC/Silicon Tubing
☐	☒	Teflon/Silicon Tubing
☐	☒	Airlift
☐	☒	Hand Pump
☐	☒	In-line Filter
☐	☒	Press/Vac Filter

(✓ All That Apply at Location)

☐ Methanol (100%)
☐ 25% Methanol/75% ASTM Type II water
☒ Deionized Water
☒ Liquinox Solution
☐ Hexane
☐ HNO₃/D.I. Water Solution
☐ Potable Water
☐ None

Field Analysis Data

PID: Ambient Air 0 ppm Well Mouth 0 ppm Purge Data Collected ☐ In-line ☒ In Container ☐ Turbid ☒ Clear ☐ Cloudy ☐ Colored ☐ Odor

Sample Observations:

Purge Data	@ <u>2</u> (11/29/94) Gal.	@ <u>4</u> Gal.	@ <u>Purged</u> Gal.	@ <u> </u> Gal.	@ <u> </u> Gal.
Temperature, Deg. C	<u>NA</u>	<u>9.9</u>	<u> </u>	<u> </u>	<u> </u>
pH, units	<u> </u>	<u>8.3</u>	<u> </u>	<u> </u>	<u> </u>
Specific Conductivity (<u>ms</u> μ mhos/cm)	<u> </u>	<u>1416</u>	<u> </u>	<u> </u>	<u> </u>
Turbidity (NTUS)	<u> </u>	<u>50</u>	<u> </u>	<u> </u>	<u> </u>
Oxidation - Reduction, +/- mv	<u> </u>	<u>NA</u>	<u> </u>	<u> </u>	<u> </u>
Dissolved Oxygen, ppm	<u> </u>	<u>8.8 NO COR</u>	<u> </u>	<u> </u>	<u> </u>
Salinity μ S/cm	<u> </u>	<u>0.06</u>	<u> </u>	<u> </u>	<u> </u>

Sample Collection Requirements
(✓ If Required at this Location)

Analytical Parameter	✓ If Sample Collected	Preservation Method	Volume Required	Sample Bottle ILot Nos.
☒ VOCs	☒	4°C	2x40 ml	<u> </u>
☒ SVOCs	☒	4°C	2x1 liter AG	<u> </u>
☒ Metals	☒	HNO ₃ , 4°C	1x1 liter P	<u> </u>
☒ Cyanide	☒	NaOH, 4°C	1x500mLP	<u> </u>
☐ Nitrate/Sulfate	☐	H ₂ SO ₄ , 4°C	1x1 liter P	<u> </u>
☐ Nitrate/Phosphate	☐	H ₂ SO ₄ , 4°C	1x1 liter P	<u> </u>
☒ Pest/PCB	☒	4°C	3x1 liter AG	<u> </u>
☐ TPH	☐	H ₂ SO ₄ , 4°C	2x1 liter AG	<u> </u>
☐ TOC	☐	H ₂ SO ₄ , 4°C	1x1 liter P	<u> </u>

Notes: visited and Purged MW-101 to dryness on 10/29/94; returned next day, purged to dryness, sampled as well recharged. collected extra volume for Duplicate (ENMW101XXXX94KD) and MS/MSD. DO meter erratic.

GROUNDWATER SAMPLE FIELD DATA RECORD

Project: ENR, INC

Site: MW-102

Project Number: 717C-40

Date: 11/30/94

Sample Location ID: ENMW1C2XXXX94XX

Time: Start: 0730 End: 0750

Signature of Sampler: Rick Buttl

Water Level/Well Data

Well Depth 14.8 Ft. ☒ Measured ☐ Historical ☒ Top of Well ☐ Top of Protective Casing

Well Riser Stick-up -0.2 Ft. (from ground) Flush Mount Protective -0.2 Ft. Casing/Well Difference

Protective 0 Ft. Casing

Depth to Water 5.69 Ft. (5.06' on 11/29/94) Well Material: ☒ PVC ☐ SS Well Locked?: ☒ Yes ☐ No Well Dia. ☒ 2 inch ☐ 4 inch ☐ 6 inch

Water Level Equip. Used: ☒ Elect. Cond. Probe ☐ Float Activated ☐ Press. Transducer

Height of Water Column 9.11 Ft. X 1.5 Gal/Ft. (2 in.) 1.5 Gal/Ft. (4 in.) 1.5 Gal/Ft. (6 in.) 1.5 Gal/Ft. (in.)

Well Integrity: Prot. Casing Secure ☐ Yes ☐ No Concrete Collar Intact ☐ Other ☐

Total Gal Purged 1.5 Gal/Vol.

Equipment Documentation

Purging/Sampling Equipment Used:

(✓ If Used For)

Purging	Sampling	Equipment ID
☒	☒	Peristaltic Pump
☐	☐	Submersible Pump
☐	☒	Bailer
☐	☒	PVC/Silicon Tubing
☐	☒	Teflon/Silicon Tubing
☐	☐	Airlift
☐	☒	Hand Pump
☐	☒	In-line Filter
☐	☐	Press/Vac Filter

Decontamination Fluids Used:

(✓ All That Apply at Location)

- ☐ Methanol (100%)
- ☐ 25% Methanol/75% ASTM Type II water
- ☐ Deionized Water
- ☐ Liquinox Solution
- ☐ Hexane
- ☐ HNO₃/D.I. Water Solution
- ☐ Potable Water
- ☐ None

Field Analysis Data

PID: Ambient Air 0 ppm Well Mouth 0 ppm Purge Data Collected ☐ In-line ☒ In Container ☐ Turbid ☒ Clear ☐ Cloudy ☐ Colored ☐ Odor

Purge Data @ 2 Gal. @ 4 Gal. @ Gal. @ Gal. @ Gal.

Parameter	2 Gal.	4 Gal.	Gal.	Gal.	Gal.
Temperature, Deg. C	<u>NA</u>	<u>9.8</u>			
pH, units	<u>5.5</u>	<u>5.62</u>			
Specific Conductivity (µmhos/cm)	<u>2.30</u>	<u>42</u>			
Turbidity (NTUS)	<u>NA</u>	<u>NA</u>			
Oxidation - Reduction, mV	<u>NA</u>	<u>NA</u>			
Dissolved Oxygen, ppm	<u>NA</u>	<u>8.63 - m. cont</u>			
Salinity (‰)	<u>NA</u>	<u>0.10</u>			

Sample Collection Requirements (✓ If Required at this Location)

Analytical Parameter	✓ If Sample Collected	Preservation Method	Volume Required	Sample Bottle ILot Nos.
☒ VOCs	☒	4°C	2x40 ml	
☒ SVOCs	☒	4°C	2x1 liter AG	
☒ Metals	☒	HNO ₃ , 4°C	1x1 liter P	
☒ Cyanide	☒	NaOH, 4°C	1x500mLP	
☐ Nitrate/Sulfate	☐	H ₂ SO ₄ , 4°C	1x1 liter P	
☐ Nitrate/Phosphate	☐	H ₂ SO ₄ , 4°C	1x1 liter P	
☒ Pest/PCB	☒	4°C	3x1 liter AG	
☐ TPH	☐	H ₂ SO ₄ , 4°C	2x1 liter AG	
☐ TOC	☐	H ₂ SO ₄ , 4°C	1x1 liter P	

Notes: visited and Purged MW-101 to dryness on 11/29/94; returned next day, purged to dryness, sampled as well recharged. DC meter readings erratic.

GROUNDWATER SAMPLE FIELD DATA RECORD

Project: ENRX, Inc.
 Project Number: 7-70-40

Site: MW-103
 Date: 11/30/94

Sample Location ID: ENMW103XXXX94XX

Time: Start: 0835 End: 1015
 Signature of Sampler: Bl But

Water Level/Well Data

Well Depth 8.42 Ft. ☒ Measured ☐ Historical ☒ Top of Well ☐ Top of Protective Casing
 Well Riser Stick-up -0.53 Ft. (from ground) Flush Mt. Protective -0.53 Ft. Casing/Well Difference
 Protective 0 Ft. Casing
 Depth to Water 2.35 Ft. Well Material: ☒ PVC ☐ SS Well Locked?: ☒ Yes ☐ No Well Dia. ☒ 2 inch ☐ 4 inch ☐ 6 inch
 Water Level Equip. Used: ☒ Elect. Cond. Probe ☐ Float Activated ☐ Press. Transducer
 Height of Water Column 6.07 Ft. ☒ .16 Gal/Ft. (2 in.) ☐ .65 Gal/Ft. (4 in.) ☐ 1.5 Gal/Ft. (6 in.) ☐ Gal/Ft. (in.)
 [1.0 Gal/Vol. 4.0 Total Gal Purged] Well Integrity: Prot. Casing Secure ☒ Yes ☐ No Concrete Collar Intact ☒ Yes ☐ No Other ☐ Yes ☐ No

Equipment Documentation

Purging/Sampling Equipment Used:

Decontamination Fluids Used:

(✓ If Used For)
 Purging ☒ Sampling ☒
 Equipment ID
 Peristaltic Pump ☐
 Submersible Pump ☐
 Bailer ☐
 PVC/Silicon Tubing ☐
 Teflon/Silicon Tubing ☐
 Airlift ☐
 Hand Pump ☐
 In-line Filter ☐
 Press/Vac Filter ☐
 (✓ All That Apply at Location)
 Methanol (100%) ☐
 25% Methanol/75% ASTM Type II water ☐
 Deionized Water ☒
 Liquinox Solution ☒
 Hexane ☐
 HNO₃/D.I. Water Solution ☐
 Potable Water ☐
 None ☐

Field Analysis Data

PID: Ambient Air 0 ppm Well Mouth 0 ppm Purge Data Collected ☒ In-line ☐ In Container Sample Observations: Turbid ☒ Clear ☐ Cloudy
 Colored ☒ Odor ☐

Purge Data	@ 1.0 Gal.	@ 2.0 Gal.	@ 3.0 Gal.	@ 4.0 Gal.	@ Gal.
Temperature, Deg. C	8.9	8.6	9.3	9.3	
pH, units	7.9	7.8	7.7	7.7	
Specific Conductivity (µmhos/cm)	2.22	1.51	1.35	1.3	
Turbidity (NTUS)	23	6	4	2	
Oxidation - Reduction, +/- mv	276	NA	NA	NA	
Dissolved Oxygen, ppm	0.10	0.95	0.7	0.5	
Salinity		0.06	0.06	0.05	

Sample Collection Requirements

(✓ If Required at this Location)

Analytical Parameter	✓ If Sample Collected	Preservation Method	Volume Required	Sample Bottle (Lot Nos.)
VOCs	☒	4°C	2x40 ml	
SVOCs	☒	4°C	2x1 liter AG	
Metals	☒	HNO ₃ , 4°C	1x1 liter P	
Cyanide	☒	NaOH, 4°C	1x500mLP	
Nitrate/Sulfate	☐	H ₂ SO ₄ , 4°C	1x1 liter P	
Nitrate/Phosphate	☐	H ₂ SO ₄ , 4°C	1x1 liter P	
Pest/PCB	☒	4°C	3x1 liter AG	
TPH	☐	H ₂ SO ₄ , 4°C	2x1 liter AG	
TOC	☐	H ₂ SO ₄ , 4°C	1x1 liter P	

Notes: water color initially black during purging - changed to clear w/ no color.

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SECTION 3.0
ANALYTICAL DATA TABLES

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ABB Environmental Services
Data Usability Report
ENRX, Inc., 7170-40
2/15/95

Introduction

This memo summarizes the usability of the analytical results generated for the ENRX, Inc. Site. Laboratory analyses were performed in accordance with the New York State Department of Environmental Conservation (NYSDEC) Analytical Services Protocol (ASP), and the data were validated using the criteria specified by U.S. Environmental Protection Agency (USEPA) Region II, modified to include NYSDCE requirements. A detailed evaluation of the laboratory quality control results (QC) is provided in the Data Validation Report.

Usability is based on validated sample results. Rejected results ("R" qualifier) represent unusable data since the analyte presence or absence is uncertain. In general, sample results with qualifiers other than "R" are considered usable. Laboratory data from the ENRX, Inc. Site will be used to determine whether hazardous wastes have been disposed at the site and to evaluate the potential threat to human health and the environment.

The data validation summary attached indicates which laboratory results are considered non-compliant when compared to the ASP requirements. However, the majority of these non-compliant results represent minor quality control problems and do not affect data usability. The cases where quality control problems affected usability and/or resulted in the rejection of data are discussed in the following sections. In most cases these problems are typical analytical difficulties or are the result of sample matrix problems.

Volatile Organics

The volatile organic compounds (VOCs) analyses were acceptable and may be considered suitable for their intended use. Methylene chloride and acetone, common laboratory contaminants, were detected in the laboratory method blank and the equipment blank. All sample results less than the action level (i.e., 10 times the blank concentration for common contaminants, 5 times for other contaminants) were reported as non-detect. Some calibration problems (continuing calibration percent differences outside acceptance limits) were observed, which represent typical laboratory performance. The affected compounds were qualified as estimated, and this minor deficiency does not affect usability. In sample ENWT105XXX94XX, positive results for trichloroethene, tetrachloroethene, acetone, carbon disulfide, 2-butanone, toluene and methylene chloride were qualified as estimated due to system monitoring compound recoveries above the acceptance range, indicating a potential high bias. 1,1-Dichloroethene; trichloroethene; and chlorobenzene did not meet RPD criteria in the MS/MSD performed on sample ENSS102XXX94XX and its duplicate. Positive and non detected results for 1,1-dichloroethene; trichloroethene; and chlorobenzene were qualified as estimated. 1,1-Dichloroethene did not meet RPD criteria in the MS/MSD performed on sample ENMW101XXX94XX and its duplicate. Positive and non detected

results for 1,1 Dichloroethene were qualified as estimated. Benzene and toluene did not meet RPD criteria in the MS/MSD performed on sample ENCD103XXX94XX and its duplicate. Positive and non detected results for benzene and toluene were qualified as estimated. These represent typical quality control problems and do not affect the useability of this data.

Positive results for vinyl chloride, acetone, trichloroethene, tetrachloroethene, toluene, ethyl benzene, and total xylenes in sample ENCL102XXX94XX and its field duplicate were qualified as estimated due to RPD criteria not being met. Positive results for 1,1,1 trichloroethane, trichloroethene, tetrachloroethane, and total xylenes for sample ENWT101XXX94XX and its field duplicate were qualified as estimated due to RPD criteria being out of control limits. Positive results for acetone and 4-methyl-2-pentanone for sample ENMW101XXX94XX and its field duplicate were qualified as estimated due to RPD criteria being out of control limits, indicating a lab precision problem. However, this does not affect the useability of this data. The internal standard area in samples ENWT103XXX94XX, ENWT104XXX94XX, ENWT105XXX94XX and ENWT106XXX94XX was out of control limits and positive and non detected results for 2-hexanone; 4-methyl-2-pentanone; tetrachloroethene; 1,1,2,2-tetrachloroethane; toluene; chlorobenzene; ethylbenzene; styrene; and total xylene were qualified as estimated. 1,4-Difluorobenzene was below acceptance criteria for sample ENWT103XXX94XX. Positive and non detected results for 1,1,1-trichloroethane; carbon tetrachloride; bromodichloromethane; 1,2-dichloropropane; trans-1,3 dichloropropene; trichloroethene; dibromochloromethane; 1,1,2-trichloroethane; benzene; cis-1,3-dichloropropene; and bromoform were qualified as estimated. This does not affect the useability of this data.

Semivolatile Organics

The semivolatile organic compounds (SVOCs) analyses provided acceptable results, and the values may be used as presented. Bis(2-ethylhexyl)phthalate, a common laboratory contaminant, was detected in the laboratory method blank. All sample results less than the action level (i.e., 10 times the blank concentration for common contaminants, 5 times for other contaminants) were reported as non-detect. Some calibration problems (continuing calibration percent differences outside acceptance limits) were observed, which represent typical laboratory performance. However, non detected results for hexachlorocyclopentadiene and 2,4-dinitrophenol were rejected due to percent difference and RRF criteria not being met. The rejected results should not be used to determine the absence of these compounds in the affected samples. Recovery was less than 10% for all compounds in the MS/MSD performed on sample ENWT101XXX94XX and its duplicate. Positive results were qualified as estimated and non detected results were rejected. Recovery was less than 10% for pentachlorophenol in the MS/MSD performed on sample ENSS102XXX94XX and its duplicate. Positive results for pentachlorophenol were qualified as estimated and non detected results were rejected. The rejected results should not be used to determine the absence of these compounds in the affected samples.

Some precision problems (Field duplicate RPD out of acceptance limits) were observed. The affected compounds were qualified as estimated, and this minor deficiency does not affect usability. Samples ENCD101XXX94XX and ENWT102XXX94XX were qualified as

estimated because of their low total solids content. This qualification does not affect the usability of these semivolatile data. A low relative response factor was observed in the calibration for perylene associated with samples ENCL104XXX94XX and ENCL103XXX94XX. Positive and non detected results were qualified as estimated for associated compounds.

Pesticides/PCBs

The pesticides/PCBs results are acceptable and may be used as presented. Samples ENCD101XXX94XX and ENWT102XXX94XX were qualified as estimated because of their low total solids content. Results for endrin aldehyde, Alpha-BHC, delta-BHC, gamma-BHC, gamma chlordane, and 4,4'-DDT were qualified as estimated because standard linearity criteria were not met, but usability of these data is not affected. Several samples exhibited surrogate recoveries below the acceptance limits, indicating a potential low bias, and the results were qualified as estimated. These low recoveries may be attributable to matrix problems and do not affect the usability of the results. However, recovery was 0% for one surrogate resulting in the rejection of non detected results in samples ENSS101XXX94XX, ENCD101XXX94XX, ENCD103XXX94XX, ENWT102XXX94XX, ENWT104XXX94XX, and ENWT105XXX94XX. The rejected data should not be used to determine the absence of these compounds in the affected samples.

Due to blank contamination, methoxychlor was qualified as non detected in associated samples where the results were below the calculated blank action level. Sample ENWT101XXX94XX was extracted 14 days beyond acceptance limits for holding times. All results for the sample were qualified as estimated. All matrix spike compounds exceeded percent recovery and RPD criteria for sample ENCL102XXX94XX and its duplicate. Positive and non detected results for gamma-BHC, heptachlor, aldrin, dieldrin, endrin, and 4,4'-DDT were qualified as estimated, indicating a potential high bias. In the MS/MSD performed on sample ENWT101XXX94XX and its duplicate, all compounds were outside percent recovery criteria. Positive results were qualified as estimated and non detected results were rejected for gamma BHC, heptachlor, aldrin, dieldrin, and 4,4'-DDT. The rejected results should not be used to determine the absence of these compounds in the affected samples. In the MS/MSD performed on sample ENWT101XXX94XX and its duplicate, aldrin, endrin and 4,4'-DDT did not meet RPD criteria. Positive and non detected results for these compounds were qualified as estimated. In the MS/MSD performed on sample ENSS102XXX94XX and its duplicate, heptachlor and dieldrin were above RPD criteria. Positive and non detected results for these compounds were qualified as estimated in this sample and its duplicate. This does not affect the useability of this data. Some precision problems (field duplicate RPD out of acceptance criteria) were observed. The affected samples were qualified as estimated. This does not affect the useability of this data. Due to percent difference criteria being out of the acceptance range, alpha chlordane was qualified as estimated for samples ENWT104XXX94XX and ENWT105XXX94XX. Gamma-BHC was rejected for samples ENCD101XXX94XX, ENWT101XXX94XX, and ENWT103XXX94XX. Heptachlor epoxide was rejected for sample ENWT101XXX94XX. 4,4'-DDT was rejected for samples ENCD103XXX94XX, ENWT103XXX94XX, ENWT104XXX94XX, ENSS102XXX94XD, ENTW105XXX94XX, and

ENWT101XXX94XX. Endrin aldehyde was rejected for samples ENCL102XXX94XX, ENWT101XXX94XX, and ENWT101XXX94XD. Gamma chlordane was rejected in sample ENWT101XXX94XX. Alpha-BHC was rejected for sample ENWT101XXX94XX. Aldrin was rejected in samples ENWT101XXX94XX and ENWT102XXX94XX. Endosulfan I was rejected for samples ENWT101XXX94XX, ENWT101XXX94XD, and ENWT104XXX94XX. Heptachlor was rejected for sample ENWT101XXX94XD. Dieldrin was rejected for sample ENWT103XXX94XX DL. Beta-BHC was rejected for sample ENWT106XXX94XX. Endosulfan II was rejected for samples ENSS101XXX94XX and ENSS102XXX94XD. The rejected results should not be used to determine the absence of these compounds in the affected samples. Additional compounds were also out of acceptance criteria resulting in the estimation of their results in associated samples. This does not affect the useability of this data.

Inorganics

The majority of the inorganics analyses are acceptable and may be used as presented. Spike recoveries for silver, cadmium, antimony, and chromium were below acceptance limits and recoveries for lead were above acceptance limits, indicating an accuracy problem. The associated results were qualified as estimated. These low recoveries may be attributable to matrix and/or laboratory problems and do not affect the usability of the results. Lead contamination above the acceptance criterion was observed in an equipment blank. Positive lead results below the action level were rejected in associated samples. Results should not be used to determine the absence of lead in site samples. Percent recovery was less than 30% for selenium in the aqueous matrix spike analysis. Results were rejected for all associated samples. The rejected results should not be used to determine the absence of selenium in these samples. Percent recovery for lead in the matrix spike was over 200% and lead results were rejected in associated samples. This should not be used to determine the absence of lead in associated samples. Percent recoveries outside of acceptance criteria were observed in other samples resulting in the estimation of their results. This represents typical laboratory performance and does not affect data useability. Duplicate injection readings for selenium and thallium were below acceptance criteria for percent recovery and results for selenium and thallium were qualified as estimated in associated samples. Potassium, cadmium, and silver did not meet %D criteria for serial dilution. All results for potassium, cadmium and silver greater than 10 times the IDL were qualified as estimated in associated samples. This does not affect data useability. Due to low solids content, positive and non detected results for samples ENCD101XXX94XX and ENWT102XXX94XX were qualified as estimated. Some precision problems were observed with field duplicates not meeting RPD criteria. Positive and non detected results were qualified as estimated in associated samples.

EP Toxicity/ Hazardous waste characteristics

Cadmium and chromium were observed in the preparation blank as negative values. Positive and non detected results less than 10 times the IDL were qualified as estimated. Percent recovery for the MS/MSD performed on barium was 150%. Positive barium results were rejected in associated samples. The rejected results should not be used to determine the

absence of barium in associated samples. Laboratory duplicate analysis was not performed on some samples and the results of these were qualified as estimated. This does not affect the useability of the data. Field duplicate criteria were not met for chromium in sample ENSS102XXX94XX and its field duplicate. Positive and non detected results for chromium were qualified as estimated in these samples. Field duplicate criteria were not met for lead in sample ENWT101XXX94XX and its duplicate. Positive and non detected results for lead were qualified as estimated in these samples.

Tentatively Identified Compounds (TICs)

The ASP analytical procedures for volatile and semivolatile organics may also detect the presence of additional compounds which are not included on the Target Compound List. The mass spectra of these non-target compounds (up to 10 VOCs and 20 SVOCs) are compared to library spectra using a computerized search routine, and the best matches are evaluated by the laboratory. If a good library match can not be made the compound is reported as "unknown". The concentrations are estimated by comparing the compound's response to the that of the closest internal standard.

Corrosivity, ignitability, reactive cyanide, reactive sulfide

All quality control criteria were met for these parameters. The results represent typical laboratory performance.

Data Quality Objectives (DQOs)

DQOs are based on the premise that different data uses require different levels of data quality. Data quality refers to the degree of uncertainty of analytical data with respect to precision, accuracy, representativeness, completeness, and comparability (PARCC). These objectives are established based on site conditions, the purpose of the field program, and the knowledge of the measurement systems used for generation of the analytical data.

No major quality control problems were observed during the data validation process which would affect the usability of the sample results. A discussion of the laboratory data quality as it relates to the PARCC objectives is presented below.

Precision and Accuracy

Precision refers to the reproducibility of a measurement under certain specified conditions, and accuracy measures the bias associated with the sampling and analysis process. Precision and accuracy are affected by both field and laboratory conditions. Precision was monitored through the analysis of field and laboratory duplicate samples; accuracy was measured through the analysis of field and laboratory blanks, matrix spikes, and surrogate spikes. The ASP protocols used for the analysis of samples define the criteria for acceptable precision and accuracy. No major precision and accuracy problems were observed which would affect usability. Some matrix spike recoveries (listed previously) were outside of acceptance criteria, indicating a potential accuracy problem. In general, these deficiencies are considered minor and do not affect useability.

Several target analytes were reported at concentrations less than the CRQL (and were qualified as estimated, "J"). Uncertainty exists for the quantitation of concentrations less than the CRQL. While these results provide information on the presence of contamination, these values should be qualified for use in decisions. In some cases precision between the two columns used for pesticides/PCBs analyses was outside the acceptance limit, and the results were qualified with a "P" by the laboratory. While these concentrations should be considered as estimated, they provide an indication of contamination and are suitable for use.

Representativeness

Measurements are made so that the results obtained are representative of the sampling population, the medium (e.g. soil and groundwater), and the site conditions. The sampling protocols were developed to ensure that the samples were representative of the media, that sampling locations were properly selected, and that a sufficient number of samples were collected. Sample handling protocols (chain-of-custody, storage, and transportation) were adequate to preserve the sample integrity. Proper documentation established that the correct protocols had been followed. Co-located samples (field duplicates) were also collected to assess representativeness, and no major problems were observed which would affect usability.

Completeness

The characteristic of completeness is defined as the percent of valid data obtained as compared to what would be expected under normal conditions. The USEPA has found that CLP protocols typically generate data that is 80% complete. Because sampling activities are often influenced by field conditions the ENRX, Inc. Site Work Plan provided estimates of the number of samples to be collected during the field program. There were no significant deviations from the proposed field program. Corrosivity, ignitability, reactive cyanide, reactive sulfide, and volatile organic analyses were found to be 100% complete. Semivolatile organic and inorganic analyses were found to be 99% complete. Pesticide analyses were 96% complete and EP toxicity analyses were found to be 93% complete.

Comparability

The characteristic of comparability reflects both the internal consistency of measurements and the expression of results in units which are consistent with other organizations reporting similar data. Each value reported for a given measurement should be similar to other values within the same data set and with other related data sets. Comparability was assured through the use of standardized sampling procedures and ASP analytical methods.



ABB Environmental Services
Data Validation Report
ENRX, Inc., 7170-40
February 15, 1995

I. INTRODUCTION

This report summarizes the data validation results for the data packages generated by Nytest Environmental Inc. concerning the water and soil samples collected from October 4 to November 30, 1994. Review was performed in accordance with the U.S. Environmental Protection Agency (USEPA) *National Functional Guidelines for Organic Data Review (June 1991)* and *Laboratory Data Validation Functional Guidelines for Evaluating Inorganics Analyses (October 1989)*, along with the appropriate USEPA Region II and New York State Department of Environmental Conservation (NYSDEC) revisions. The data tables referred to in this memo consist of the following:

Table 1: Laboratory Report of Analysis

Table 2: Validation / Summary Table

Table 1 presents the analytical results as reported by the laboratory. Table 2 presents the validated results with the appropriate data qualifiers. The laboratory qualifiers used on Table 1 are defined in Attachment I; data validation qualifiers used on Table 2 are defined in Attachment II. For all organics analyses, compound results below the contract required quantitation limit (CRQL) were flagged with a J by the laboratory on Table 1. These results were considered estimated and flagged J on Table 2. Compound results greater than the calibration range were flagged with an E by the laboratory and on Table 1. Samples containing these compounds were diluted and reanalyzed, and the diluted results flagged with a D by the laboratory and on Table 1. On Table 2, the diluted results for all compounds beyond calibration range were inserted into the original results and the remainder of the diluted analysis deleted from Table 2. Pesticides/PCBs that had greater than 25% difference between the two analytical columns were flagged with a P by the laboratory. Pesticides/PCBs that had greater than 25% but less than 50% difference between the two analytical columns were considered estimated and were flagged J on Table 2. When the difference between the two analytical columns was greater than 50% but less than 90% the pesticide/PCB compound was considered tentatively identified, and the associated result represents an estimated concentration (JN). If the difference between the two analytical columns was greater than 90% the pesticide/PCB result was rejected (R). For all inorganic analyses, analyte results below the contract required detection limit (CRDL) were flagged with a B by the laboratory on Table 1. These results were considered estimated and were flagged with a J on Table 2.

The samples were analyzed using the following methods:

- Target Compound List (TCL) Volatile Organic Compounds (VOCs) - NYSDEC Analytical Services Protocol (ASP) 91-1
- TCL Semivolatile Organic Compounds (SVOCs) - NYSDEC ASP 91-2
- TCL Pesticides/PCBs - NYSDEC ASP 91-3
- TCL Inorganics - NYSDEC ASP Contract Laboratory Program Superfund Methods
- EP Toxicity metal, ignitability, corrosivity, and reactivity, USEPA SW-846

This narrative presents a summary of the laboratory quality control (QC) deficiencies and the resulting qualification of the data.

II. VOLATILE ORGANIC COMPOUNDS

A. Holding Times

Holding times are evaluated to address the validity of the results based on the elapsed time from Validated Time of Sample Receipt (VTSR) to analysis. All samples were analyzed within the required 7 day holding time.

B. Gas Chromatograph/Mass Spectrometer (GC/MS) Instrument Performance Check

Bromofluorobenzene (BFB) is analyzed every 12 hours to verify the instrument's mass resolution, identification, and sensitivity. All BFB ion abundance criteria were met.

C. GC/MS Initial and Continuing Calibration

Initial calibration demonstrates instrument linearity and ensures that the instrument can produce acceptable qualitative and quantitative results. The initial calibration relative standard deviation (RSD) must be less than 30%, and the relative response factor (RRF) must be greater than 0.05. Continuing calibration checks are performed every 12-hours to demonstrate that the instrument can produce acceptable qualitative and quantitative results as established by the initial calibration. The continuing calibration percent difference (%D) must be less than 25%, and the RRF must be greater than 0.05. Chloromethane, methylene chloride, acetone, 2-butanone, and 2-hexanone did not meet RSD criteria for initial calibration standards. Methylene chloride, acetone and 2-butanone positive results were qualified as estimated for associated samples. No action was required for chloromethane or 2-hexanone. Chloroethane, acetone, and 2-hexanone did not meet percent difference criteria for continuing calibration standards. Positive acetone results were estimated in the associated samples. No action was required for chloroethane or 2-hexanone.

D. Blanks

Laboratory (method) and field (trip/equipment) blanks are analyzed to determine the presence and magnitude of contamination resulting from field or laboratory activities. Action levels are

calculated at 5 times the concentration in the associated blank (10 times for methylene chloride, acetone, and 2-butanone). Sample results below this action level are considered attributable to blank contamination; results greater than this level are considered to be acceptable. Due to trip, equipment, field, or laboratory method blank contamination, methylene chloride, acetone, and toluene were qualified as non-detect in associated samples where the results were below the calculated blank action level.

E. System Monitoring Compounds Recoveries

System monitoring compounds are added to all samples and blanks prior to analyses to assess recovery (accuracy). Action is taken if any system monitoring compound recoveries are outside the acceptance range, or if any one recovery is less than 10%. System monitoring compound recoveries for ENWT105XXX94XX were above the method acceptance range, indicating a potential high bias. The positive results (i.e., trichloroethene, tetrachloroethene, acetone, carbon disulfide, 2-butanone, toluene and methylene chloride) for this sample were qualified as estimated.

F. Matrix Spike/Matrix Spike Duplicate (MS/MSD)

MS/MSD analyses are performed at a frequency of 5% to assess method precision and accuracy. Action is taken if recoveries are outside the acceptance range, or if the relative percent difference (RPD) for spiked compounds is above the control limit. A total of five MS/MSD analyses were performed, two for low level aqueous samples, two for low level soil samples, and one for medium level soil samples. 1,1-Dichloroethene did not meet RPD acceptance limits in the MS/MSD performed on aqueous sample ENMW101XXX94XX. Positive and non-detected 1,1-dichloroethene results in sample ENMW101XXX94XX and its field duplicate ENMW101XXX94XX were qualified as estimated. Toluene and chlorobenzene did not meet percent recovery criteria, and benzene and toluene did not meet RPD criteria in the MS/MSD performed on soil sample ENCD103XXX94XX. Positive and non-detected benzene and toluene results for sample ENCD103XXX94XX were qualified as estimated. No action was required for chlorobenzene. 1,1-Dichloroethene did not meet percent recovery criteria or RPD criteria in the MS/MSD performed on soil sample ENSS102XXX94XX. Trichloroethene and chlorobenzene did not meet RPD criteria in the same MS/MSD. Positive and non-detected 1,1-dichloroethene; trichloroethene; and chlorobenzene results were qualified as estimated for sample ENSS102XXX94XX and its field duplicate ENSS102XXX94XD.

G. Field Duplicates

Field duplicate samples are collected and analyzed to assess sampling and analytical precision. Field duplicate control limits are: RPD of less than 30% for water samples and 50% for soil samples. Vinyl chloride, acetone, trichloroethene, tetrachloroethene, toluene, ethylbenzene, and total xylenes did not meet the control limit in aqueous sample ENCL102XXX94XX and its duplicate, ENCL102XXX94XD. Positive results for those compounds for both samples were qualified as estimated. 1,1,1-Trichloroethane; trichloroethene; tetrachloroethane; and total

xylenes did not meet the control limit in soil sample ENWT101XXX94XX and its duplicate, ENWT101XXX94XD. Positive results for those compounds for both samples were qualified as estimated. Acetone and 4-methyl-2-pentanone did not meet the control limit in aqueous sample ENMW101XXX94XX and its duplicate, ENMW101XXX94XD. Positive results for those compounds for sample ENMW101XXX94XX were qualified as estimated. No action was required for ENMW101XXX94XD because acetone and 4-methyl-2-pentanone were non-detected.

H. Internal Standard Response

The internal standard response is monitored for each sample to verify GC/MS sensitivity and the stability of the detector's response. The internal standard area must be >50% and <100%, and the retention time must be within $\pm 30\%$, of the associated calibration standard. Chlorobenzene-d5 was below acceptance criteria for the following samples: ENWT103XXX94XX, ENWT104XXX94XX, ENWT105XXX94XX, and ENWT106XXX94XX. Positive and non-detected results for all compounds for which chlorobenzene-d5 is the internal standard (i.e., 2-hexanone; 4-methyl-2-pentanone; tetrachloroethene; 1,1,2,2-tetrachloroethane; toluene; chlorobenzene; ethylbenzene; styrene; and total xylene) were qualified as estimated in the samples listed. 1,4-Difluorobenzene was below acceptance criteria for ENWT103XXX94XX. Positive and non-detected results for all compounds for which 1,4-Difluorobenzene is the internal standard (i.e., 1,1,1-trichloroethane; carbon tetrachloride; bromodichloromethane; 1,2-dichloropropane; trans-1,3-dichloropropene; trichloroethene; dibromochloromethane; 1,1,2-trichloroethane; benzene; cis-1,3-dichloropropene; and bromoform) were qualified as estimated in the sample.

I. Target Compound Identification

Chromatograms and mass spectra are reviewed to minimize the reporting of false positive and false negatives. For each compound detected, the relative retention time must be within ± 0.06 units and the qualitative criteria for mass spectral identification must be met. No problems were observed.

J. Compound Quantitation

Laboratory calculations were checked to verify that reported concentrations and CRQLs were accurate. The calculations which were reviewed were performed correctly, and the CRQLs were adjusted for sample size and dilution factors. Soil sample percent solid content is evaluated to determine whether the sample was correctly classified as a soil. If solid content falls between 10% and 50% positive and non-detected results are estimated. If solid content is less than 10% results are calculated and reported as an aqueous sample. No sample results were qualified due to low solid content.

K. Tentatively Identified Compounds (TICs)

All TIC spectra were reviewed to verify that the identifications were acceptable, laboratory contamination was taken into account, and the correct assignments of compound classes were made. Reported concentrations represent estimated (J) values.

III. SEMIVOLATILE ORGANIC COMPOUNDS

A. Holding Times

Holding times are evaluated to address the validity of the results based on the elapsed time from VTSR to analysis. All samples were analyzed within the required 5-day holding time for extraction and the 40-day holding time from extraction to analysis.

B. GC/MS Instrument Performance Check

Decafluorotriphenylphosphine (DFTPP) is analyzed every 12 hours to verify the instrument's mass resolution, identification, and sensitivity. All DFTPP ion abundance criteria were met.

C. GC/MS Initial and Continuing Calibration

Initial calibration demonstrates instrument linearity and ensures that the instrument can produce acceptable qualitative and quantitative results. The initial calibration RSD must be less than 30%, and the RRF must be greater than 0.05. Continuing calibration checks are performed every 12-hours to demonstrate that the instrument can produce acceptable qualitative and quantitative results as established by the initial calibration. The continuing calibration %D must be less than 25%, and the RRF must be greater than 0.05. Bis(2-chloroethyl)ether, hexachlorocyclopentadiene, diethylphthalate, 4-chlorophenyl-phenylether, and benzo(k)fluoranthene exceed 30% RSD, but are below 50% RSD. No action was required for these compounds. 3-Nitroaniline and hexachlorocyclopentadiene exceed 50% RSD. Associated positive and non-detected sample results are qualified as estimated. Hexachlorocyclopentadiene and 2,4-dinitrophenol did not meet percent difference or RRF criteria for continuing calibration standards. Positive results for those compounds were qualified as estimated and non-detected results were rejected in associated samples. Bis(2-chloroethyl)ether; N-nitroso-di-n-propylamine; 2-nitrophenol; 4-chloroaniline; hexachlorocyclopentadiene; 3-nitroaniline; 2,4-dinitrophenol; 4-nitrophenol; diethylphthalate; 4-chlorophenyl-phenylether; 4-nitroaniline; 4,6-dinitro-2-methylphenol; pentachlorophenol; 3,3'-dichlorobenzidine; di-n-octylphthalate; and benzo(k)fluoranthene exceed 25% difference, but are below 50% difference. Positive results for those compounds were qualified as estimated for associated samples. 2,4-Dinitrophenol; 4-nitroaniline; 4,6-dinitro-2-methylphenol; N-nitrosodiphenylamine; and 3,3'-dichlorobenzidine exceeded 50% difference. Positive and non-detected results were qualified as estimated for associated samples.

D. Blanks

Laboratory (method) and field (equipment) blanks are analyzed to determine the presence and magnitude of contamination resulting from field or laboratory activities. Action levels are calculated at 5 times the concentration in the associated blank (10 times for phthalates). Sample results below this action level are considered attributable to blank contamination; results greater than this level are considered to be acceptable. Due to equipment, field, or laboratory method blank contamination, bis(2-ethylhexyl)phthalate was qualified as non-detect in associated samples where the results were below the calculated blank action level.

E. Surrogate Recoveries

Surrogate compounds are added to all samples and blanks prior to extraction to assess recovery (accuracy). Action is taken if any two acid or base/neutral surrogates are outside the acceptance range, or if any one surrogate recovery is less than 10%. All surrogate recoveries were within QC limits.

F. Matrix Spike/Matrix Spike Duplicate

MS/MSD analyses are performed at a frequency of 5% to assess method precision and accuracy. Action is taken if recoveries are outside the acceptance range, or if the RPD for spiked compounds is above the control limit. A total of four MS/MSD analyses were performed: two for aqueous samples and two for soil samples. Recovery was below 10% for all compounds in the MS/MSD performed on soil sample ENWT101XXX94XX. Positive results for all matrix spike compounds were qualified as estimated, and non-detected results were rejected for sample ENWT101XXX94XX and its field duplicate ENWT101XXX94XD. Percent recoveries for pentachlorophenol and pyrene in the MS/MSD performed on aqueous sample ENCL102XXX94XX were above QC limits. Positive pyrene results were qualified as estimated for sample ENCL102XXX94XD. No action was required for pentachlorophenol in either the original sample or its field duplicate, or for pyrene in the original sample. Recovery was below 10% for pentachlorophenol in the MS performed on soil sample ENSS102XXX94XX. Pentachlorophenol did not meet percent recovery criteria in the MSD performed on soil sample ENSS102XXX94XX, or RPD criteria between the MS and MSD concentrations. Positive results for pentachlorophenol were qualified as estimated, and non-detected results were rejected for sample ENSS102XXX94XX and its field duplicate ENSS102XXX94XD. 1,4-dichlorobenzene RPD was above QC limits for the MS/MSD performed on soil sample ENSS102XXX94XX. Positive and non-detected 1,4-dichlorobenzene results were qualified as estimated for sample ENSS102XXX94XX and its field duplicate ENSS102XXX94XD.

G. Field Duplicates

Field duplicate samples are collected and analyzed to assess sampling and analytical precision. Field duplicate control limits are: RPD of less than 30% for water samples and 50% for soil

samples. Pyrene and di-n-octylphthalate did not meet the control limit in aqueous sample ENCL102XXX94XX and its duplicate, ENCL102XXX94XD. Positive results for those compounds for both samples were qualified as estimated. Phenanthrene, fluoranthene, pyrene, and chrysene did not meet the control limit for soil sample ENWT101XXX94XX and its duplicate, ENWT101XXX94XD. Positive results for those compounds for both samples were qualified as estimated. Phenanthrene, fluoanthene, pyrene, benzo(a)anthracene, chrysene, benzo(b)fluoranthene, benzo(k)fluoranthene, benzo(a)pyrene, indeno(1,2,3-cd)pyrene, and benzo(g,h,i)perylene did not meet the control limit for soil sample ENSS102XXX94XX and its field duplicate ENSS102XXX94XD. Positive results for those compounds for both samples were qualified as estimated.

H. Internal Standard Response

The internal standard response is monitored for each sample to verify GC/MS sensitivity and the stability of the detector's response. The internal standard area must be >50% and <100%, and the retention time must be within $\pm 30\%$, of the associated calibration standard. Perylene-d12 response was low, but greater than 25% of associated continuing calibration standard response, for samples ENCL104XXX94XX DL and ENCL103XXX94XX. Positive and non-detected results for all compounds for which perylene-d12 is the internal standard (i.e., di-n-octylphthalate, benzo(b)fluoranthene, benzo(k)fluoranthene, benzo(a)pyrene, indeno(1,2,3-cd)pyrene, dibenz(a,h)anthracene, and benzo(g,h,i)perylene) were qualified as estimated in the samples listed.

I. Target Compound Identification

Chromatograms and mass spectra are reviewed to minimize the reporting of false positive and false negatives. For each compound detected, the relative retention time must be within ± 0.06 units and the qualitative criteria for mass spectral identification must be met. No problems were observed.

J. Compound Quantitation

Laboratory calculations were checked to verify that reported concentrations and CRQLs were accurate. The calculations which were reviewed were performed correctly, and the CRQLs were adjusted for sample size and dilution factors. Soil sample percent solid content is evaluated to determine whether the sample was correctly classified as a soil. If solid content falls between 10% and 50% positive and non-detected results are estimated. If solid content is less than 10% results are calculated and reported as an aqueous sample. Positive and non-detected results for samples ENCD101XXX94XX and ENWT102XXX94XX were qualified as estimated due to low solid content.

K. Tentatively Identified Compounds

All TIC spectra were reviewed to verify that the identifications were acceptable, laboratory contamination was taken into account, and the correct assignments of compound classes were made.. Reported concentrations represent estimated (J) values.

IV. PESTICIDES/PCBs

A. Holding Times

Holding times are evaluated to address the validity of the results based on the elapsed time from VTSR to analysis. Sample extraction must be performed within 5 days of VTSR, and sample analysis must be done within 40 days of extraction. Sample ENWT101XXX94XD was extracted 14 days beyond the required holding time; all results for the sample were qualified as estimated.

B. Instrument Performance Check

Performance checks are performed to verify target compound resolution and the instrument's sensitivity. All compound resolution criteria were met for the resolution check mixture and the performance evaluation mixture (PEM). For the PEM, retention time criteria were met, the %D between the calculate and true concentration was less than 25%, and the endrin and 4,4'-DDT breakdown was less than 20% (less than 30% for endrin and 4,4'-DDT combined).

C. Initial and Continuing Calibration

Initial calibration demonstrates instrument linearity and ensures that the instrument can produce acceptable qualitative and quantitative results. For the initial calibration linearity criteria to be met, the RSD must be less than 20%. Standards were run at the required frequency. Alpha-BHC, delta-BHC, gamma-BHC, endrin aldehyde, gamma-chlordane, and 4,4'-DDT did not meet the RSD criterion. Positive and non-detected results for alpha-BHC, delta-BHC, gamma-BHC, endrin aldehyde, gamma-chlordane, and 4,4'-DDT were qualified as estimated in associated samples. Continuing calibration checks are performed to demonstrate that the instrument can produce acceptable qualitative and quantitative results as established by the initial calibration. The continuing calibration %D must be less than 25%, and the retention times must fall within the windows established by the initial calibration. Standards were run at the required frequency, and these criteria were met.

D. Blanks

Laboratory (method) and field (equipment/source) blanks are analyzed to determine the presence and magnitude of contamination resulting from field or laboratory activities. Action levels are calculated at 5 times the concentration in the associated blank. Sample results below this action level are considered attributable to blank contamination; results greater than this level are

considered to be acceptable. Due to equipment, field, or laboratory method blank contamination, methoxychlor was qualified as non-detect in associated samples where the results were below the calculated blank action level.

E. Surrogate Recoveries

Surrogate compounds are added to all samples and blanks prior to extraction to assess recovery (accuracy). Action is taken if both surrogates are outside the acceptance range in either column, or if any one surrogate recovery is less than 10%. Recovery was 0% for one surrogate in samples ENSS101XXX94XX, ENCD101XXX94XX, ENCD103XXX94XX, ENWT102XXX94XX, ENWT104XXX94XX, and ENWT105XXX94XX. Positive results were qualified as estimated and non-detected results were rejected in these samples due to surrogate percent recovery. Surrogate percent recoveries did not meet acceptance criteria in samples ENWT101XXX94XD or ENMW103XXX94XX. Positive results (i.e., 4,4'-DDT and endrin aldehyde in sample ENWT101XXX94XD, none in sample ENMW103XXX94XX) obtained from the column that did not meet surrogate percent recovery criteria, and all non-detected results in those samples were qualified as estimated.

F. Matrix Spike/Matrix Spike Duplicate

Matrix spike/matrix spike duplicate analyses are performed at a frequency of 5% to assess method precision and accuracy. Action is taken if recoveries are outside the acceptance range, or if the relative percent difference (RPD) for spiked compounds (RSD for unspiked compounds) is above the control limit. A total of four MS/MSD analyses were performed: two for aqueous samples and two for soil samples. All matrix spike compounds exceeded percent recovery and RPD acceptance limits in the MS/MSD performed on aqueous sample ENCL102XXX94XX. Positive and non-detected gamma-BHC (Lindane), heptachlor, aldrin, dieldrin, endrin, and 4,4'-DDT results were qualified as estimated for the sample ENCL102XXX94XX and its field duplicate ENCL102XXX94XD. All matrix spike compounds were outside of percent recovery criteria in the MS/MSD performed on soil sample ENWT101XXX94XX. Positive results were qualified as estimated and non-detected results were rejected for gamma-BHC (Lindane), heptachlor, aldrin, dieldrin, and 4,4'-DDT in the sample and its field duplicate ENWT101XXX94XD. No action was necessary for endrin. Aldrin, endrin and 4,4'-DDT did not meet RPD criteria in the MS/MSD performed on soil sample ENWT101XXX94XX. Positive and non-detected results for aldrin, endrin and 4,4'-DDT were qualified as estimated for sample ENWT101XXX94XX and its field duplicate ENWT101XXX94XD. Heptachlor and dieldrin were above RPD criteria in the MS/MSD performed on soil sample ENSS102XXX94XX. Positive and non-detected results for heptachlor and dieldrin in sample ENSS102XXX94XX and its field duplicate ENSS102XXX94XD were qualified as estimated.

G. Field Duplicates

Field duplicate samples are collected and analyzed to assess sampling and analytical precision. Field duplicate control limits are: RPD of less than 30% for water samples and 50% for soil samples. Alpha-BHC, gamma-BHC, heptachlor, aldrin, heptachlor epoxide, endosulfan I, 4,4'-DDE, endrin aldehyde, 4,4'-DDT, and gamma chlordane did not meet the control limit in sample ENWT101XXX94XX and its duplicate, ENWT101XXX94XD. Positive results for those compounds for both samples were qualified as estimated.

H. Cleanup Checks

Cleanup procedures (gel permeation chromatography (GPC) and florisil) are used to remove interferences from sample extracts. Cleanup checks verify acceptable recovery of pesticides through the cleanup process. Recoveries must be between 80 - 120% (florisil) and 80 - 110% (GPC). These criteria were met.

I. Target Compound Identification

Chromatograms and mass spectra are reviewed to minimize the reporting of false positive and false negatives. For each compound detected, the retention time must be within the retention time window determined during initial calibration for both the primary and confirmation columns, and the percent difference between the results obtained from each column must be less than 25%. If the percent difference between the two is between 25% and 50% the compound result is qualified as estimated. If the percent difference between the two is between 50% and 90% the compound result is qualified as an analyte that is tentatively identified and whose associated result represents an estimated concentration. If the percent difference between the two is greater than 90% the compound result is rejected. Alpha chlordane was qualified as estimated for samples ENWT104XXX94XX and ENWT105XXX94XX. Gamma-BHC (Lindane) was rejected for sample ENCD101XXX94XX, ENWT101XXX94XX, and ENWT103XXX94XX. Heptachlor epoxide was rejected for sample ENWT101XXX94XX, and qualified as JN for sample ENWT105XXX94XX. 4,4'-DDD was rejected for samples ENCD103XXX94XX, ENWT103XXX94XX, and ENWT104XXX94XX. Endrin aldehyde was rejected for samples ENCL102XXX94XX, ENWT101XXX94XD, and ENWT101XXX94XX. 4,4'-DDE was qualified as estimated for samples ENCL104XXX94XX and ENWT105XXX94XX, qualified as JN for sample ENWT104XXX94XX, and rejected for samples ENWT101XXX94XX and ENWT103XXX94XX. Gamma chlordane was rejected for sample ENWT101XXX94XX, and qualified as estimated for sample ENWT103XXX94XX. Alpha-BHC was qualified as JN for sample ENCD101XXX94XX, and rejected for sample ENWT101XXX94XX. Aldrin was rejected for samples ENWT101XXX94XX and ENWT102XXX94XX, and qualified as estimated for sample ENWT103XXX94XX. Endosulfan I was rejected for samples ENWT101XXX94XX, ENWT101XXX94XD, and ENWT104XXX94XX, and qualified as JN for sample ENTWT105XXX94XX. Heptachlor was rejected for sample ENWT101XXX94XD and qualified as JN for sample ENTWT105XXX94XX. 4,4'-DDT was qualified as estimated for sample

ENWT101XXX94XD, as JN for samples ENSS101XXX94XX and ENWT103XXX94XX, and rejected for samples ENSS102XXX94XD and ENTWT105XXX94XX. Dieldrin was rejected for sample ENWT103XXX94XX, qualified as JN for sample ENWT104XXX94XX, and as estimated for sample ENTWT105XXX94XX. Beta-BHC was rejected for sample ENWT106XXX94XX DL. Endosulfan II was rejected for samples ENSS101XXX94XX and ENSS102XXX94XD. Methoxychlor was qualified as estimated for sample ENSS102XXX94XX. Endrin ketone was rejected for samples ENSS102XXX94XX and ENSS102XXX94XD. Arochlor 1260 was qualified as estimated in sample ENBS102XX694XX.

J. Compound Quantitation

Laboratory calculations were checked to verify that reported concentrations and CRQLs were accurate. The calculations which were reviewed were performed correctly, and the CRQLs were adjusted for sample size and dilution factors. Soil sample percent solid content is evaluated to determine whether the sample was correctly classified as a soil. If solid content falls between 10% and 50% positive and non-detected results are estimated. If solid content is less than 10% results are calculated and reported as an aqueous sample. Positive and non-detected results for samples ENCD101XXX94XX and ENWT102XXX94XX were qualified as estimated due to low solid content.

V. INORGANICS

A. Holding Times

Holding times are evaluated to address the validity of the results based on the elapsed time from VTSR to analysis. All samples were analyzed within the following holding times:

metals - 6 months
mercury - 26 days
cyanide - 12 days

B. Calibration

Calibration demonstrates instrument linearity and ensures that the instrument can produce acceptable qualitative and quantitative results. The minimum number of standards were analyzed by the laboratory as specified by the method. For the initial calibration linearity criteria to be met, the correlation coefficient must be greater than 0.995 for furnace atomic absorption (AA), mercury, and cyanide. Initial calibration verification (ICV) and continuing calibration verification (CCV) results must be between 90 -110% (80 - 120% for mercury and 85 - 115% for cyanide). For the CRDL standard, recoveries must be between 80 -120%. All standards were run at the required frequency. Positive and non-detected cadmium, antimony, silver, and chromium results within the affected range (i.e., standard true value ± 2 times the CRDL) were estimated for associated water samples because the CRDL standard recovery was below the acceptance limits.

Positive lead results within the affected range (i.e., standard true value ± 2 times the CRDL) were estimated for associated water samples because the CRDL standard recovery was above the acceptance limits. Positive and non-detected cadmium, antimony, and silver results within the affected range (i.e., standard true value ± 2 times the CRDL) were estimated for associated soil samples because the CRDL standard recovery was below the acceptance limits. Positive lead results within the affected range (i.e., standard true value ± 2 times the CRDL) were estimated for associated soil samples because the CRDL standard recovery was above the acceptance limits.

C. Blanks

Laboratory (preparation/calibration) and field (equipment/source) blanks are analyzed to determine the presence and magnitude of contamination resulting from field or laboratory activities. If contamination above the CRDL is found in a field, equipment, or method blank an action level is calculated. The action level is calculated at 5 times the concentration in the blank. Positive sample results below this action level are rejected; results greater than this level are considered to be acceptable. Lead contamination above the acceptance criterion was found in an equipment blank. Positive lead results below the action level were rejected in associated soil samples because of this blank contamination.

D. Interference Check Sample (ICS)

The ICS verifies the instrument's interelement and background correction factors for inductively coupled plasma (ICP) analyses. The ICS recovery must be within 80 - 120%. All results were reviewed and found to be acceptable.

E. Laboratory Control Sample (LCS)

The LCS monitors the overall laboratory performance from sample preparation through analysis. Aqueous LCS recoveries must fall between 80 - 120%, and solid LCS results must fall within the limits established by USEPA for that LCS. All results were reviewed and found to be acceptable.

F. Laboratory Duplicate Analysis

Duplicate results provide a measure of the laboratory's analytical precision. The RPD must be less than 50% (100% for soil) for sample results ≥ 5 times the CRDL, or $\pm$ CRDL (± 2 times the CRDL for soil) for sample results less < 5 times the CRDL. All duplicate results were within acceptance limits. All duplicate results were within acceptance limits.

G. Matrix Spike

Matrix spike analyses are performed to assess method accuracy. Spike recoveries must fall within the range of 75 - 125%. Four matrix spike analyses were performed: two for aqueous samples and two for soil samples. Percent recovery was less than 30% for selenium in aqueous matrix

spike analysis. Results were rejected for all associated aqueous samples. Manganese percent recovery was above the acceptance limit for aqueous matrix spike analysis. Positive results were qualified as estimated for associated aqueous samples. Percent recoveries were below acceptance limits for antimony, arsenic, selenium, silver, and thallium in one soil matrix spike analysis. Positive and non-detected results for those analytes were qualified as estimated in associated samples. Percent recoveries were above acceptance limits for arsenic, lead, selenium, vanadium, and cyanide in one soil matrix spike analysis. No action was required for selenium or cyanide. Positive arsenic and vanadium results were qualified as estimated in associated samples. Positive lead results were rejected in associated samples because matrix spike percent recovery was above 200%.

H. Furnace AA QC

Duplicate injections and post digestion spikes provide a measure of precision and accuracy for furnace AA analyses. Duplicate injections must be within 20% RSD, and spike recoveries must be between 85 - 115%. All RSDs were reviewed and found to be acceptable. Percent recoveries were low for selenium in samples ENCL102XXX94XD, ENCL102XXX94XX, ENCL103XXX94XX, ENCL104XXX94XX, ENWT106XXX94XX, ENBS101XX694XX, ENSS101XXX94XX, and ENMW102XXX94XX, and for thallium in samples ENCL103XXX94XX, ENCL104XXX94XX, ENWT101XXX94XD, ENWT101XXX94XX, ENBS101XX694XX, ENSS101XXX94XX. Positive and non-detected results for these analytes in these samples were qualified as estimated. Method of standard additions (MSA) is performed for sample quantitation if upon analysis of the sample and its analytical spike the sample absorbance or concentration is greater than or equal to 50% of the spike and the spike recovery is less than 85% or greater than 115%. MSA is evaluated for degree of dependence between concentration and absorbance in the concentration range of the MSA standards. Correlation coefficient must be greater than 0.995. Quality control criteria were met for all MSAs performed.

I. ICP Serial Dilution

Serial dilution analyses evaluate the effects of physical or chemical interferences in the sample matrix. Serial dilution results must agree within 10%D of the original sample for initial concentration greater than 10 times the IDL. Potassium and cadmium aqueous serial dilution results did not meet quality control criteria. All potassium and cadmium results greater than 10 times the IDL were qualified as estimated in associated samples. Cadmium and silver soil serial dilution results did not meet quality control criteria. All cadmium and silver results greater than 10 times the IDL were qualified as estimated in associated samples.

J. Sample Result Verification

Laboratory calculations were checked to verify that reported concentrations and IDLs were accurate. The calculations which were reviewed were performed correctly, and the CRDLs were

adjusted for sample size and dilution factors. Soil sample percent solid content is evaluated to determine whether the sample was correctly classified as a soil. If solid content falls between 10% and 50% positive and non-detected results are estimated. If solid content is less than 10% results are calculated and reported as an aqueous sample. Positive and non-detected results for samples ENCD101XXX94XX and ENWT102XXX94XX were qualified as estimated due to low solid content.

K. Field Duplicates

Field duplicate samples are collected and analyzed to assess sampling and analytical precision. Field duplicate results must be within the following criteria: difference of less than CRDL for water (2 times the CRDL for soil) when results are <5 times the CRDL or RPD <50% for water samples (100% for soil samples) when results are ≥ 5 times the CRDL). Aqueous sample ENCL102XXX94XX and its duplicate ENCL102XXX94XD did not meet quality control criteria for nickel. Positive and non-detected nickel results were qualified as estimated in the sample and its duplicate only. Aqueous sample ENMW101XXX94XX and its duplicate ENMW101XXX94XD did not meet quality control criteria for aluminum or lead. Positive and non-detected results for those analytes were qualified as estimated in the sample and its duplicate only. Soil sample ENWT101XXX94XX and its duplicate ENWT101XXX94XD did not meet quality control criteria for sodium, vanadium, or cyanide. Positive and non-detected results for those analytes were qualified as estimated in the sample and its duplicate only.

V. EP TOXICITY METALS

EP toxicity metals analyses were evaluated for hold times, calibration, blank contamination, interference check sample, laboratory control sample, laboratory duplicate analysis, matrix spike, furnace AA quality control (when applicable), ICP serial dilution, sample result verification, and field duplicate. Preparation blanks had negative results for cadmium and chromium. Positive and non-detected sample results less than 10 times the IDL were qualified as estimated. Laboratory duplicate analysis was not performed for batches associated with the following samples: ENSS101XXX94XX, ENSS102XXX94XX, and ENSS102XXX94XD. Positive results were qualified as estimated. Laboratory duplicate analysis was not performed for lead by furnace AA for the following samples: ENCL103XXX94XX and ENCL104XXX94XX. Positive results were qualified as estimated. Laboratory duplicate results were not within acceptance limits (acceptance limits are similar to those for laboratory duplicate results in CLP inorganics analysis, but IDL is used instead of CRDL in the criteria) for chromium, lead (by ICP), or silver. Positive and non-detected chromium, lead, and silver results in all associated samples were qualified as estimated. Matrix spike analysis was not performed for batches associated with the following samples: ENSS101XXX94XX, ENSS102XXX94XX, and ENSS102XXX94XD. Positive cadmium, chromium and lead results less than 4 times the spiking level were qualified as estimated. Matrix spike analysis exceeded percent recovery criteria (criteria are similar to those for matrix spike recovery in CLP inorganics analysis) for barium. Positive barium results were rejected in associated samples because percent recovery was greater than 150%. Furnace AA

analytical spike percent recovery criteria (criteria are similar to those for analytical spike recovery in CLP inorganics analysis) were not met for selenium in samples ENCL103XXX94XX and ENCL104XXX94XX, or for arsenic in sample ENCL104XXX94XX. Positive and non-detected results for these analytes in these samples were qualified as estimated. Field duplicate criteria (criteria are similar to those for field duplicate results in CLP inorganics analysis, but IDL is used instead of CRDL in the criteria) were not met for chromium in sample ENSS102XXX94XX and its field duplicate ENSS102XXX94XD, or for lead in sample ENWT101XXX94XX and its field duplicate ENWT101XXX94XD. Positive and non-detected results for these analytes in the associated samples were qualified as estimated.

VI. CORROSIVITY, IGNITABILITY, REACTIVE CYANIDE, REACTIVE SULFIDE

Corrosivity, ignitability, reactive cyanide and reactive sulfide analyses were evaluated for hold times, calibration, method blank contamination (there is no method blank for corrosivity or ignitability), laboratory control sample, matrix spike, and field duplicate. All quality control criteria were met.

**Attachment I - Definition of Laboratory Qualifiers
(for Table 1 - Laboratory Report of Analysis)**

Organic Data Qualifiers

- J - Indicates an estimated concentration below the contract required detection level (CRQL) but greater than 0 or when estimating a concentration for TICs.
- U - Indicates that compound was analyzed but not detected. The sample quantitation limit is adjusted for dilution and percent moisture.
- B - Indicates analyte was detected in both the sample and the associated laboratory method blank.
- E - Indicates that the analyte concentration exceeded the calibration range of the GC/MS and that a re-analysis of a diluted sample is required.
- D - Indicates that sample concentration was obtained by dilution to bring result within calibration range.
- N - Indicates presumptive evidence of a compound. This flag is used for TICs where the identification is based on a library search and is applied to all TIC results. For general classes of compounds (hydrocarbons, etc.) this flag is not used.
- P - This flag is used for pesticides/PCBs when there is greater than 25% difference between the concentrations on the two columns used for analysis. The lower value is reported.
- C - This flag applies to pesticide/PCBs results when the identification has been confirmed by GC/MS.
- A - Indicates that a TIC is a suspected aldol-condensation product.
- X - Laboratory-defined qualifier used to provide additional information not covered by the other qualifiers.

Inorganic Data Qualifiers

- E - The reported concentration is estimated because of the presence of an interference.
- M - Duplicate injection precision criteria were not met.
- N - Spiked sample recovery not within control limits.
- s - The reported concentration was determined by the method of standard additions.
- W - Post-digestion spike for furnace atomic adsorption analysis is outside control limits.
- B - Concentration reported is below CRDL but greater than the IDL.
- * - Duplicate analysis not within control limits.
- + - Correlation coefficient for the method of standard additions was less than 0.995
- U - Indicates that compound was analyzed but not detected. The sample quantitation limit is adjusted for dilution and percent moisture.

**Attachment II - Definition of Validation Qualifiers
(for Table 2 - Validation/Summary Table)**

- J - Estimated concentration because QC criteria were not met.
- R - Results were rejected because of serious QC deficiencies.
- U - Indicates that compound was analyzed but not detected. The sample quantitation limit is adjusted for dilution and percent moisture.
- N - Indicates presumptive evidence of a compound. This flag is used for TICs where the identification is based on a library search and is applied to all TIC results. For general classes of compounds (hydrocarbons, etc.) this flag is not used.
- UJ - Quantitation limit was estimated concentration because QC criteria were not met.
- JN - Presence of an analyte was tentatively identified and the associated result represents an estimated concentration.

Table 1
Laboratory Report of Analysis

ANALYTE	LOCATION:			
	ISIS ID:			
	LAB NUMBER:			
	DATE SAMPLED:			
	DATE ANALYZED:			
	QS-003	QS-004	QS-XX1	QS-2
	ENQS003XXX94XX	ENQS004XXX94XX	ENQSXX1XXX94XX	ENQSXX2XXX94XX
	2241108	2263807	2223319	2223313
	10/28/94	11/30/94	10/04/94	10/04/94
	11/08/94	12/02/94	10/13/94	10/13/94
	SOW-3/90 - 11	CRQL		
Chloromethane	10	10 U	10 U	10 U
Bromomethane	10	10 U	10 U	10 U
Vinyl Chloride	10	10 U	10 U	10 U
Chloroethane	10	10 U	10 U	10 U
Methylene Chloride	10	4 JB	10 U	15 B
Acetone	10	25	8 J	10 U
Carbon Disulfide	10	10 U	10 U	10 U
1,1-Dichloroethene	10	10 U	10 U	10 U
1,1-Dichloroethane	10	10 U	10 U	10 U
1,2-Dichloroethene (total)	10	10 U	10 U	10 U
Chloroform	10	10 U	10 U	10 U
1,2-Dichloroethane	10	10 U	10 U	10 U
2-Butanone	10	10 U	10 U	10 U
1,1,1-Trichloroethane	10	10 U	10 U	10 U
Carbon Tetrachloride	10	10 U	10 U	10 U
Bromodichloromethane	10	10 U	10 U	10 U
1,2-Dichloropropane	10	10 U	10 U	10 U
cis-1,3-Dichloropropene	10	10 U	10 U	10 U
Trichloroethene	10	10 U	10 U	1 J
Dibromochloromethane	10	10 U	10 U	10 U
1,1,2-Trichloroethane	10	10 U	10 U	10 U
Benzene	10	10 U	10 U	10 U
trans-1,3-Dichloropropene	10	10 U	10 U	10 U
Bromoform	10	10 U	10 U	10 U
4-Methyl-2-Pentanone	10	10 U	10 U	10 U
2-Hexanone	10	10 U	10 U	10 U
Tetrachloroethene	10	10 U	10 U	10 U
1,1,2,2-Tetrachloroethane	10	10 U	10 U	10 U
Toluene	10	2 J	10 U	2 J
Chlorobenzene	10	10 U	10 U	10 U
Ethylbenzene	10	10 U	10 U	10 U
Styrene	10	10 U	10 U	10 U
Total Xylenes	10	10 U	10 U	10 U
=====				
Dilution Factor:	1.00	1.00	1.00	1.00
Sample Volume\Weight (ml\g):	5.00	5.00	5.00	5.00
Associated Method Blank:	M0746.D	N0496.D	N9594.D	N9594.D
Associated Equipment Blank:	-	-	-	-
Associated Field Blank:	-	-	-	-
Associated Trip Blank:	-	-	-	-

Site: EQUIPMENT RINSATE

U: not detected J: estimated B: blank contamination

Table 1
Laboratory Report of Analysis

LOCATION:	QT-001	QT-002
ISIS ID:	ENQT001XXX94XX	ENQT002XXX94XX
LAB NUMBER:	2223320	2263808
DATE SAMPLED:	10/04/94	11/30/94
DATE ANALYZED:	10/13/94	12/02/94

ANALYTE	SOW-3/90 - 11	CRQL		
Chloromethane	10	10 U	10 U	
Bromomethane	10	10 U	10 U	
Vinyl Chloride	10	10 U	10 U	
Chloroethane	10	10 U	10 U	
Methylene Chloride	10	17 B	2 JB	
Acetone	10	10 U	68	
Carbon Disulfide	10	10 U	10 U	
1,1-Dichloroethene	10	10 U	10 U	
1,1-Dichloroethane	10	10 U	10 U	
1,2-Dichloroethene (total)	10	10 U	10 U	
Chloroform	10	10 U	10 U	
1,2-Dichloroethane	10	10 U	10 U	
2-Butanone	10	10 U	10 U	
1,1,1-Trichloroethane	10	10 U	10 U	
Carbon Tetrachloride	10	10 U	10 U	
Bromodichloromethane	10	10 U	10 U	
1,2-Dichloropropane	10	10 U	10 U	
cis-1,3-Dichloropropene	10	10 U	10 U	
Trichloroethene	10	10 U	10 U	
Dibromochloromethane	10	10 U	10 U	
1,1,2-Trichloroethane	10	10 U	10 U	
Benzene	10	10 U	10 U	
trans-1,3-Dichloropropene	10	10 U	10 U	
Bromoform	10	10 U	10 U	
4-Methyl-2-Pentanone	10	10 U	10 U	
2-Hexanone	10	10 U	10 U	
Tetrachloroethene	10	10 U	10 U	
1,1,2,2-Tetrachloroethane	10	10 U	10 U	
Toluene	10	10 U	10 U	
Chlorobenzene	10	10 U	10 U	
Ethylbenzene	10	10 U	10 U	
Styrene	10	10 U	10 U	
Total Xylenes	10	10 U	10 U	

Dilution Factor:	1.00	1.00
Sample Volume/Weight (ml/g):	5.00	5.00
Associated Method Blank:	N9594.D	N0496.D
Associated Equipment Blank:	-	-
Associated Field Blank:	-	-
Associated Trip Blank:	-	-

Site: TRIP BLANKS

U: not detected B: blank contamination J: estimated

Table 1
Laboratory Report of Analysis

	LOCATION:	CL-102 DUP	CL-102 DUP	CL-102	CL-102	CL-103	CL-104	CL-104
	ISIS ID:	ENCL102XXX94XD	ENCL102XXX94XD	ENCL102XXX94XX	ENCL102XXX94XX	ENCL103XXX94XX	ENCL104XXX94XX	ENCL104XXX94XX
	LAB NUMBER:	2223317	2223317 D	2223314	2223314 D	2223318	2223312	2223312 D
	DATE SAMPLED:	10/04/94	10/04/94	10/04/94	10/04/94	10/04/94	10/05/94	10/05/94
	DATE ANALYZED:	10/13/94	10/14/94	10/13/94	10/14/94	10/13/94	10/13/94	10/13/94
ANALYTE	SOW-3/90 - 11	CRQL						
Chloromethane	10	10 U	100 U	10 U	100 U	10 U	200 U	10000 U
Bromomethane	10	10 U	100 U	10 U	100 U	10 U	200 U	10000 U
Vinyl Chloride	10	34	68 JD	80	100 U	10 U	230	10000 U
Chloroethane	10	10 U	100 U	10 U	100 U	10 U	200 U	10000 U
Methylene Chloride	10	18 B	120 BD	17 B	120 BD	16 B	830 B	4100 JBD
Acetone	10	16	100 U	10 U	100 U	10 U	200 U	10000 U
Carbon Disulfide	10	10 U	100 U	10 U	100 U	10 U	200 U	10000 U
1,1-Dichloroethene	10	10 U	100 U	10 U	100 U	10 U	200 U	10000 U
1,1-Dichloroethane	10	16	15 JD	15	16 JD	10 U	14000 E	12000 D
1,2-Dichloroethene (total)	10	450 E	480 D	420 E	480 D	10 U	17000 E	13000 D
Chloroform	10	10 U	100 U	10 U	100 U	10 U	200 U	10000 U
1,2-Dichloroethane	10	10 U	100 U	10 U	100 U	10 U	200 U	10000 U
2-Butanone	10	10 U	100 U	10 U	100 U	10 U	200 U	10000 U
1,1,1-Trichloroethane	10	190	190 D	170	170 D	10 U	60000 E	57000 D
Carbon Tetrachloride	10	10 U	100 U	10 U	100 U	10 U	200 U	10000 U
Bromodichloromethane	10	10 U	100 U	10 U	100 U	10 U	200 U	10000 U
1,2-Dichloropropane	10	10 U	100 U	10 U	100 U	10 U	200 U	10000 U
cis-1,3-Dichloropropene	10	10 U	100 U	10 U	100 U	10 U	200 U	10000 U
Trichloroethene	10	10	100 U	6 J	100 U	10 U	34000 E	33000 D
Dibromochloromethane	10	10 U	100 U	10 U	100 U	10 U	200 U	10000 U
1,1,2-Trichloroethane	10	10 U	100 U	10 U	100 U	10 U	200 U	10000 U
Benzene	10	10 U	100 U	10 U	100 U	10 U	200 U	10000 U
trans-1,3-Dichloropropene	10	10 U	100 U	10 U	100 U	10 U	200 U	10000 U
Bromoform	10	10 U	100 U	10 U	100 U	10 U	200 U	10000 U
4-Methyl-2-Pentanone	10	10 U	100 U	10 U	100 U	10 U	200 U	10000 U
2-Hexanone	10	10 U	100 U	10 U	100 U	10 U	200 U	10000 U
Tetrachloroethene	10	4 J	100 U	2 J	100 U	10 U	6500 E	8000 JD
1,1,2,2-Tetrachloroethane	10	10 U	100 U	10 U	100 U	10 U	200 U	10000 U
Toluene	10	28	17 JD	13	16 JD	10 U	510	10000 U
Chlorobenzene	10	10 U	100 U	10 U	100 U	10 U	200 U	10000 U
Ethylbenzene	10	8 J	100 U	3 J	100 U	10 U	43 J	10000 U
Styrene	10	10 U	100 U	10 U	100 U	10 U	200 U	10000 U
Total Xylenes	10	30	12 JD	10	100 U	10 U	210	10000 U
=====								
	Dilution Factor:	1.00	10.0	1.00	10.0	1.00	20.0	1000
	Sample Volume\Weight (ml\g):	5.00	5.00	5.00	5.00	5.00	5.00	5.00
	Associated Method Blank:	N9594.D	N9613.D	N9594.D	N9613.D	N9594.D	N9594.D	N9594.D
	Associated Equipment Blank:	ENQ5XX2XXX94XX	ENQ5XX2XXX94XX	ENQ5XX2XXX94XX	ENQ5XX2XXX94XX	ENQ5XX2XXX94XX	ENQ5XX2XXX94XX	ENQ5XX2XXX94XX
	Associated Field Blank:	-	-	-	-	-	-	-
	Associated Trip Blank:	ENQT001XXX94XX	ENQT001XXX94XX	ENQT001XXX94XX	ENQT001XXX94XX	ENQT001XXX94XX	ENQT001XXX94XX	ENQT001XXX94XX

Site: SUMP LIQUIDS

U: not detected B: blank contamination J: estimated

D: diluted result E: exceeds calibration range

Table 1
Laboratory Report of Analysis

LOCATION:	CD-101	CD-103
ISIS ID:	ENCD101XXX94XX	ENCD103XXX94XX
LAB NUMBER:	2223303	2223304
DATE SAMPLED:	10/04/94	10/04/94
DATE ANALYZED:	10/13/94	10/13/94

ANALYTE	SOW-3/90 - 11	CRQL		
Chloromethane	10	160	U	18 U
Bromomethane	10	160	U	18 U
Vinyl Chloride	10	390		18 U
Chloroethane	10	160	U	18 U
Methylene Chloride	10	280	B	32 B
Acetone	10	960		150
Carbon Disulfide	10	160	U	18 U
1,1-Dichloroethene	10	160	U	18 U
1,1-Dichloroethane	10	450		18 U
1,2-Dichloroethene (total)	10	180		18 U
Chloroform	10	160	U	18 U
1,2-Dichloroethane	10	160	U	18 U
2-Butanone	10	470		42
1,1,1-Trichloroethane	10	160		18 U
Carbon Tetrachloride	10	160	U	18 U
Bromodichloromethane	10	160	U	18 U
1,2-Dichloropropane	10	160	U	18 U
cis-1,3-Dichloropropene	10	160	U	18 U
Trichloroethene	10	120	J	18 U
Dibromochloromethane	10	160	U	18 U
1,1,2-Trichloroethane	10	160	U	18 U
Benzene	10	160	U	18 U
trans-1,3-Dichloropropene	10	160	U	18 U
Bromoform	10	160	U	18 U
4-Methyl-2-Pentanone	10	160	U	18 U
2-Hexanone	10	160	U	18 U
Tetrachloroethene	10	490		3 J
1,1,2,2-Tetrachloroethane	10	160	U	18 U
Toluene	10	260		18 U
Chlorobenzene	10	160	U	18 U
Ethylbenzene	10	380		18 U
Styrene	10	160	U	18 U
Total Xylenes	10	3800		18 U

Dilution Factor:	5.00	1.00
Percent Solids:	32	55
Sample Volume\Weight (ml\g):	5.00	5.00
Associated Method Blank:	P1171.D	P1171.D
Associated Equipment Blank:	ENQSXX1XXX94XX	ENQSXX1XXX94XX
Associated Field Blank:	-	-
Associated Trip Blank:	-	-

Site: SUMP SEDIMENTS

U: not detected B: blank contamination J: estimated

Table 1
Laboratory Report of Analysis

LOCATION:	WT-101 DUP	WT-101	WT-101
ISIS ID:	ENWT101XXX94XD	ENWT101XXX94XX	ENWT101XXX94XX
LAB NUMBER:	2223309	2223306	2223306 D
DATE SAMPLED:	10/04/94	10/04/94	10/04/94
DATE ANALYZED:	10/12/94	10/12/94	10/13/94

ANALYTE	SOW-3/90 - II	CRQL				
Chloromethane	1200		1400	U	1400	U
Bromomethane	1200		1400	U	1400	U
Vinyl Chloride	1200		1400	U	1400	U
Chloroethane	1200		1400	U	1400	U
Methylene Chloride	1200		1000	JB	980	JB
Acetone	1200		1400	U	1400	U
Carbon Disulfide	1200		1400	U	1400	U
1,1-Dichloroethene	1200		1400	U	1400	U
1,1-Dichloroethane	1200		1400	U	1400	U
1,2-Dichloroethene (total)	1200		1400	U	1400	U
Chloroform	1200		1400	U	1400	U
1,2-Dichloroethane	1200		1400	U	1400	U
2-Butanone	1200		1400	U	1400	U
1,1,1-Trichloroethane	1200		1500		7600	
Carbon Tetrachloride	1200		1400	U	1400	U
Bromodichloromethane	1200		1400	U	1400	U
1,2-Dichloropropane	1200		1400	U	1400	U
cis-1,3-Dichloropropene	1200		1400	U	1400	U
Trichloroethene	1200		610	J	1500	
Dibromochloromethane	1200		1400	U	1400	U
1,1,2-Trichloroethane	1200		1400	U	1400	U
Benzene	1200		1400	U	1400	U
trans-1,3-Dichloropropene	1200		1400	U	1400	U
Bromoform	1200		1400	U	1400	U
4-Methyl-2-Pentanone	1200		1400	U	1400	U
2-Hexanone	1200		980	J	1400	U
Tetrachloroethene	1200		10000		36000	E
1,1,2,2-Tetrachloroethane	1200		1400	U	1400	U
Toluene	1200		1400	U	1400	U
Chlorobenzene	1200		1400	U	1400	U
Ethylbenzene	1200		1400	U	1400	U
Styrene	1200		1400	U	1400	U
Total Xylenes	1200		1400	U	520	J

Dilution Factor:	1.00	1.00	5.00
Percent Solids:	84	88	
Sample Volume\Weight (ml\g):	4.00	4.00	4.00

Associated Method Blank:	N9570.D	N9570.D	N9595.D
Associated Equipment Blank:	ENQSXX1XXX94XX	ENQSXX1XXX94XX	ENQSXX1XXX94XX
Associated Field Blank:	-	-	-
Associated Trip Blank:	-	-	-

Site: WASTE-SLUDGE

U: not detected J: estimated B: blank contamination

E: exceeds calibration range D: diluted result

Table 1
Laboratory Report of Analysis

LOCATION:		WT-102	WT-103	WT-103	WT-104	WT-104	WT-105	WT-105	WT-106
ISIS ID:		ENWT102XXX94XX	ENWT103XXX94XX	ENWT103XXX94XX	ENWT104XXX94XX	ENWT104XXX94XX	ENWT105XXX94XX	ENWT105XXX94XX	ENWT106XXX94XX
LAB NUMBER:		2223305	2223310	2223310 R	2223311	2223311 R	2223301	2223301 R	2223302
DATE SAMPLED:		10/04/94	10/04/94	10/04/94	10/04/94	10/04/94	10/05/94	10/05/94	10/05/94
DATE ANALYZED:		10/12/94	10/12/94	10/13/94	10/12/94	10/13/94	10/12/94	10/12/94	10/12/94
ANALYTE	SOW-3/90 - 11	CRQL							
Chloromethane	10	20 U	17 U	17 U	17 U	17 U	11 U	11 U	11 U
Bromomethane	10	20 U	17 U	17 U	17 U	17 U	11 U	11 U	11 U
Vinyl Chloride	10	20 U	17 U	17 U	17 U	17 U	11 U	11 U	11 U
Chloroethane	10	20 U	17 U	17 U	17 U	17 U	11 U	11 U	11 U
Methylene Chloride	10	11 JB	13 JB	40 B	11 JB	63 B	11 B	12 B	11 B
Acetone	10	20 U	190	16 J	17 U	17 U	6 J	17	11 U
Carbon Disulfide	10	20 U	4 J	4 J	17 U	17 U	2 J	5 J	11 U
1,1-Dichloroethene	10	20 U	17 U	17 U	17 U	17 U	11 U	11 U	11 U
1,1-Dichloroethane	10	26	17 U	17 U	17 U	17 U	11 U	11 U	11 U
1,2-Dichloroethene (total)	10	29	17 U	17 U	17 U	17 U	11 U	11 U	11 U
Chloroform	10	20 U	17 U	17 U	17 U	17 U	11 U	11 U	11 U
1,2-Dichloroethane	10	20 U	17 U	17 U	17 U	17 U	11 U	11 U	11 U
2-Butanone	10	20 U	14 J	17 U	17 U	17 U	5 J	16	11 U
1,1,1-Trichloroethane	10	75	13 J	17 U	17 U	17 U	11 U	2 J	3 J
Carbon Tetrachloride	10	9 J	17 U	17 U	17 U	17 U	11 U	11 U	11 U
Bromodichloromethane	10	20 U	17 U	17 U	17 U	17 U	11 U	11 U	11 U
1,2-Dichloropropane	10	20 U	17 U	17 U	17 U	17 U	11 U	11 U	11 U
cis-1,3-Dichloropropene	10	20 U	17 U	17 U	17 U	17 U	11 U	11 U	11 U
Trichloroethene	10	100	14 J	16 J	6 J	17 U	23	30	7 J
Dibromochloromethane	10	20 U	17 U	17 U	17 U	17 U	11 U	11 U	11 U
1,1,2-Trichloroethane	10	20 U	17 U	17 U	17 U	17 U	11 U	11 U	11 U
Benzene	10	20 U	17 U	17 U	17 U	17 U	11 U	11 U	11 U
trans-1,3-Dichloropropene	10	20 U	17 U	17 U	17 U	17 U	11 U	11 U	11 U
Bromoform	10	20 U	17 U	17 U	17 U	17 U	11 U	11 U	11 U
4-Methyl-2-Pentanone	10	20 U	17 U	17 U	17 U	17 U	11 U	11 U	11 U
2-Hexanone	10	20 U	17 U	17 U	17 U	17 U	11 U	11 U	11 U
Tetrachloroethene	10	48	10 J	16 J	37	11 J	43	62	31
1,1,2,2-Tetrachloroethane	10	20 U	17 U	17 U	17 U	17 U	11 U	11 U	11 U
Toluene	10	20 U	2 J	2 J	17 U	17 U	4 J	7 J	5 J
Chlorobenzene	10	20 U	17 U	17 U	17 U	17 U	11 U	11 U	11 U
Ethylbenzene	10	20 U	17 U	17 U	17 U	17 U	11 U	11 U	11 U
Styrene	10	20 U	17 U	17 U	17 U	17 U	11 U	11 U	11 U
Total Xylenes	10	20 U	17 U	17 U	17 U	17 U	11 U	11 U	11 U
=====									
Dilution Factor:	1.00	1.00	1.00	1.00	1.00	1.00	1.00	1.00	1.00
Percent Solids:	49	59	59	60	60	91	91	91	
Sample Volume\Weight (ml\g):	5.00	5.00	5.00	5.00	5.00	5.00	5.00	5.00	5.00
=====									
Associated Method Blank:	P1148.D	P1148.D	P1171.D	P1148.D	P1171.D	P1148.D	P1148.D	P1148.D	P1148.D
Associated Equipment Blank:	ENQSXX1XXX94XX	ENQSXX1XXX94XX	ENQSXX1XXX94XX	ENQSXX1XXX94XX	ENQSXX1XXX94XX	ENQSXX1XXX94XX	ENQSXX1XXX94XX	ENQSXX1XXX94XX	ENQSXX1XXX94XX
Associated Field Blank:	-	-	-	-	-	-	-	-	-
Associated Trip Blank:	-	-	-	-	-	-	-	-	-

Site: WASTE-SLUDGE

U: not detected J: estimated B: blank contamination

Table 1
Laboratory Report of Analysis

LOCATION: WT-106
ISIS ID: ENWT106XXX94XX
LAB NUMBER: 2223302 R
DATE SAMPLED: 10/05/94
DATE ANALYZED: 10/12/94

ANALYTE	SOW-3/90 - 11	CRQL
Chloromethane	10	11 U
Bromomethane	10	11 U
Vinyl Chloride	10	11 U
Chloroethane	10	11 U
Methylene Chloride	10	13 B
Acetone	10	10 J
Carbon Disulfide	10	11 U
1,1-Dichloroethene	10	11 U
1,1-Dichloroethane	10	11 U
1,2-Dichloroethene (total)	10	11 U
Chloroform	10	11 U
1,2-Dichloroethane	10	11 U
2-Butanone	10	8 J
1,1,1-Trichloroethane	10	6 J
Carbon Tetrachloride	10	11 U
Bromodichloromethane	10	11 U
1,2-Dichloropropane	10	11 U
cis-1,3-Dichloropropene	10	11 U
Trichloroethene	10	12
Dibromochloromethane	10	11 U
1,1,2-Trichloroethane	10	11 U
Benzene	10	11 U
trans-1,3-Dichloropropene	10	11 U
Bromoform	10	11 U
4-Methyl-2-Pentanone	10	11 U
2-Hexanone	10	11 U
Tetrachloroethene	10	58
1,1,2,2-Tetrachloroethane	10	11 U
Toluene	10	11
Chlorobenzene	10	11 U
Ethylbenzene	10	4 J
Styrene	10	11 U
Total Xylenes	10	3 J

Dilution Factor: 1.00
Percent Solids: 91
Sample Volume\Weight (ml\g): 5.00

Associated Method Blank: P1148.D
Associated Equipment Blank: ENQSXX1XXX94XX
Associated Field Blank: -
Associated Trip Blank: -

Site: WASTE-SLUDGE

U: not detected J: estimated B: blank contamination

Table 1
Laboratory Report of Analysis

LOCATION:	SS-101	SS-102 DUP	SS-102
ISIS ID:	ENSS101XXX94XX	ENSS102XXX94XD	ENSS102XXX94XX
LAB NUMBER:	2241101	2241105	2241102
DATE SAMPLED:	10/28/94	10/28/94	10/28/94
DATE ANALYZED:	11/03/94	11/03/94	11/03/94

ANALYTE	SOW-3/90 - 11	CRQL			
Chloromethane	10		11 U	14 U	14 U
Bromomethane	10		11 U	14 U	14 U
Vinyl Chloride	10		11 U	14 U	14 U
Chloroethane	10		11 U	14 U	14 U
Methylene Chloride	10	8 JB	30 B	10 JB	
Acetone	10		11 U	14 U	14 U
Carbon Disulfide	10		11 U	14 U	14 U
1,1-Dichloroethene	10		11 U	14 U	14 U
1,1-Dichloroethane	10		11 U	14 U	14 U
1,2-Dichloroethene (total)	10		11 U	14 U	14 U
Chloroform	10		11 U	14 U	14 U
1,2-Dichloroethane	10		11 U	14 U	14 U
2-Butanone	10		11 U	14 U	14 U
1,1,1-Trichloroethane	10		11 U	14 U	14 U
Carbon Tetrachloride	10		11 U	14 U	14 U
Bromodichloromethane	10		11 U	14 U	14 U
1,2-Dichloropropane	10		11 U	14 U	14 U
cis-1,3-Dichloropropene	10		11 U	14 U	14 U
Trichloroethene	10		11 U	14 U	14 U
Dibromochloromethane	10		11 U	14 U	14 U
1,1,2-Trichloroethane	10		11 U	14 U	14 U
Benzene	10		11 U	14 U	14 U
trans-1,3-Dichloropropene	10		11 U	14 U	14 U
Bromoform	10		11 U	14 U	14 U
4-Methyl-2-Pentanone	10		11 U	14 U	14 U
2-Hexanone	10		11 U	14 U	14 U
Tetrachloroethene	10		11 U	14 U	14 U
1,1,2,2-Tetrachloroethane	10		11 U	14 U	14 U
Toluene	10		11 U	14 U	14 U
Chlorobenzene	10		11 U	14 U	14 U
Ethylbenzene	10		11 U	14 U	14 U
Styrene	10		11 U	14 U	14 U
Total Xylenes	10		11 U	14 U	14 U

Dilution Factor:	1.00	1.00	1.00
Percent Solids:	90	74	74
Sample Volume\Weight (ml\g):	5.00	5.00	5.00
Associated Method Blank:	P1621.D	P1621.D	P1621.D
Associated Equipment Blank:	ENQS003XXX94XX	ENQS003XXX94XX	ENQS003XXX94XX
Associated Field Blank:	-	-	-
Associated Trip Blank:	-	-	-

Site: SURFACE SOILS

U: not detected J: estimated B: blank contamination

Table 1
Laboratory Report of Analysis

LOCATION:	BS-101	BS-102
ISIS ID:	ENBS101XX694XX	ENBS102XX694XX
LAB NUMBER:	2241106	2241107
DATE SAMPLED:	10/27/94	10/28/94
DATE ANALYZED:	11/03/94	11/03/94

ANALYTE	SOW-3/90 - II	CRQL		
Chloromethane	10	13	U	12 U
Bromomethane	10	13	U	12 U
Vinyl Chloride	10	13	U	12 U
Chloroethane	10	13	U	12 U
Methylene Chloride	10	24	B	21 B
Acetone	10	21	B	5 JB
Carbon Disulfide	10	13	U	12 U
1,1-Dichloroethene	10	13	U	12 U
1,1-Dichloroethane	10	13	U	12 U
1,2-Dichloroethene (total)	10	13	U	12 U
Chloroform	10	13	U	12 U
1,2-Dichloroethane	10	13	U	12 U
2-Butanone	10	4	J	12 U
1,1,1-Trichloroethane	10	13	U	12 U
Carbon Tetrachloride	10	13	U	12 U
Bromodichloromethane	10	13	U	12 U
1,2-Dichloropropane	10	13	U	12 U
cis-1,3-Dichloropropene	10	13	U	12 U
Trichloroethene	10	13	U	12 U
Dibromochloromethane	10	13	U	12 U
1,1,2-Trichloroethane	10	13	U	12 U
Benzene	10	13	U	12 U
trans-1,3-Dichloropropene	10	13	U	12 U
Bromoform	10	13	U	12 U
4-Methyl-2-Pentanone	10	13	U	12 U
2-Hexanone	10	13	U	12 U
Tetrachloroethene	10	13	U	12 U
1,1,2,2-Tetrachloroethane	10	13	U	12 U
Toluene	10	13	U	12 U
Chlorobenzene	10	13	U	12 U
Ethylbenzene	10	13	U	12 U
Styrene	10	13	U	12 U
Total Xylenes	10	13	U	12 U

Dilution Factor:	1.00	1.00
Percent Solids:	77	81
Sample Volume\Weight (ml\g):	5.00	5.00
Associated Method Blank:	P1621.D	P1621.D
Associated Equipment Blank:	ENQS003XXX94XX	ENQS003XXX94XX
Associated Field Blank:	-	-
Associated Trip Blank:	-	-

Site: SOIL BORINGS

U: not detected B: blank contamination J: estimated

Table 1
Laboratory Report of Analysis

LOCATION:	MW-101 DUP	MW-101	MW-102	MW-103
ISIS ID:	ENMW101XXX94XD	ENMW101XXX94XX	ENMW102XXX94XX	ENMW103XXX94XX
LAB NUMBER:	2263804	2263801	2263805	2263806
DATE SAMPLED:	11/30/94	11/30/94	11/30/94	11/30/94
DATE ANALYZED:	12/02/94	12/05/94	12/02/94	12/02/94

ANALYTE	SOW-3/90 - 11	CRQL				
Chloromethane	10		10 U	10 U	10 U	10 U
Bromomethane	10		10 U	10 U	10 U	10 U
Vinyl Chloride	10		10 U	10 U	10 U	10 U
Chloroethane	10		10 U	10 U	10 U	6 J
Methylene Chloride	10	6 JB	8 JB	10 U	10 U	10 U
Acetone	10	10 U	44	11	34	
Carbon Disulfide	10	10 U	10 U	10 U	10 U	10 U
1,1-Dichloroethene	10	10 U	10 U	10 U	10 U	10 U
1,1-Dichloroethane	10	10 U	10 U	10 U	10 U	2 J
1,2-Dichloroethene (total)	10	10 U	10 U	10 U	10 U	10 U
Chloroform	10	10 U	10 U	10 U	10 U	10 U
1,2-Dichloroethane	10	10 U	10 U	10 U	10 U	10 U
2-Butanone	10	10 U	10 U	10 U	10 U	8 J
1,1,1-Trichloroethane	10	10 U	10 U	10 U	10 U	10 U
Carbon Tetrachloride	10	10 U	10 U	10 U	10 U	10 U
Bromodichloromethane	10	10 U	10 U	10 U	10 U	10 U
1,2-Dichloropropane	10	10 U	10 U	10 U	10 U	10 U
cis-1,3-Dichloropropene	10	10 U	10 U	10 U	10 U	10 U
Trichloroethene	10	10 U	10 U	10 U	10 U	10 U
Dibromochloromethane	10	10 U	10 U	10 U	10 U	10 U
1,1,2-Trichloroethane	10	10 U	10 U	10 U	10 U	10 U
Benzene	10	10 U	10 U	10 U	10 U	10 U
trans-1,3-Dichloropropene	10	10 U	10 U	10 U	10 U	10 U
Bromoform	10	10 U	10 U	10 U	10 U	10 U
4-Methyl-2-Pentanone	10	10 U	5 J	10 U	10 U	10 U
2-Hexanone	10	10 U	10 U	10 U	10 U	10 U
Tetrachloroethene	10	10 U	10 U	10 U	10 U	10 U
1,1,2,2-Tetrachloroethane	10	10 U	10 U	10 U	10 U	10 U
Toluene	10	10 U	10 U	10 U	10 U	10 U
Chlorobenzene	10	10 U	10 U	10 U	10 U	10 U
Ethylbenzene	10	10 U	10 U	10 U	10 U	10 U
Styrene	10	10 U	10 U	10 U	10 U	10 U
Total Xylenes	10	10 U	10 U	10 U	10 U	1 J

Dilution Factor:	1.00	1.00	1.00	1.00
Sample Volume\Weight (ml\g):	5.00	5.00	5.00	5.00
Associated Method Blank:	N0496.D	N0519.D	N0496.D	N0496.D
Associated Equipment Blank:	ENQS004XXX94XX	ENQS004XXX94XX	ENQS004XXX94XX	ENQS004XXX94XX
Associated Field Blank:	-	-	-	-
Associated Trip Blank:	ENQT002XXX94XX	ENQT002XXX94XX	ENQT002XXX94XX	ENQT002XXX94XX

Site: MONITORING WELLS

U: not detected J: estimated B: blank contamination

Table 1
Laboratory Report of Analysis

LOCATION:	QS-003	QS-004	QS-XX1	QS-XX1
ISIS ID:	ENQS003XXX94XX	ENQS004XXX94XX	ENQSXX1XXX94XX	ENQSXX1XXX94XX
LAB NUMBER:	2241108	2263807	2223319	2223319 R
DATE SAMPLED:	10/28/94	11/30/94	10/04/94	10/04/94
DATE EXTRACTED:	11/02/94	12/06/94	10/11/94	10/11/94
DATE ANALYZED:	12/02/94	12/22/94	11/18/94	11/18/94

ANALYTE	SOW-3/90 - II	CRQL			
Phenol	10	10 U	27	10 U	10 U
bis(2-Chloroethyl)ether	10	10 U	10 U	10 U	10 U
2-Chlorophenol	10	10 U	10 U	10 U	10 U
1,3-Dichlorobenzene	10	10 U	10 U	10 U	10 U
1,4-Dichlorobenzene	10	10 U	10 U	10 U	10 U
1,2-Dichlorobenzene	10	10 U	10 U	10 U	10 U
2-Methylphenol	10	10 U	10 U	10 U	10 U
2,2'-oxybis(1-Chloropropane)	10	10 U	10 U	10 U	10 U
4-Methylphenol	10	10 U	10 U	10 U	10 U
N-Nitroso-di-n-propylamine	10	10 U	10 U	10 U	10 U
Hexachloroethane	10	10 U	10 U	10 U	10 U
Nitrobenzene	10	10 U	10 U	10 U	10 U
Isophorone	10	10 U	10 U	10 U	10 U
2-Nitrophenol	10	10 U	10 U	10 U	10 U
2,4-Dimethylphenol	10	10 U	10 U	10 U	10 U
bis(2-Chloroethoxy)methane	10	10 U	10 U	10 U	10 U
2,4-Dichlorophenol	10	10 U	10 U	10 U	10 U
1,2,4-Trichlorobenzene	10	10 U	10 U	10 U	10 U
Naphthalene	10	10 U	10 U	10 U	10 U
4-Chloroaniline	10	10 U	10 U	10 U	10 U
Hexachlorobutadiene	10	10 U	10 U	10 U	10 U
4-Chloro-3-Methylphenol	10	10 U	10 U	10 U	10 U
2-Methylnaphthalene	10	10 U	10 U	10 U	10 U
Hexachlorocyclopentadiene	10	10 U	10 U	10 U	10 U
2,4,6-Trichlorophenol	10	10 U	10 U	10 U	10 U
2,4,5-Trichlorophenol	25	25 U	25 U	25 U	25 U
2-Chloronaphthalene	10	10 U	10 U	10 U	10 U
2-Nitroaniline	25	25 U	25 U	25 U	25 U
Dimethylphthalate	10	10 U	10 U	10 U	10 U
Acenaphthylene	10	10 U	10 U	10 U	10 U
2,6-Dinitrotoluene	10	10 U	10 U	10 U	10 U

Site: EQUIPMENT RINSATE

U: not detected J: estimated B: blank contamination

Table 1
Laboratory Report of Analysis

LOCATION:	QS-003	QS-004	QS-XX1	QS-XX1
ISIS ID:	ENQS003XXX94XX	ENQS004XXX94XX	ENQSXX1XXX94XX	ENQSXX1XXX94XX
LAB NUMBER:	2241108	2263807	2223319	2223319 R
DATE SAMPLED:	10/28/94	11/30/94	10/04/94	10/04/94
DATE EXTRACTED:	11/02/94	12/06/94	10/11/94	10/11/94
DATE ANALYZED:	12/02/94	12/22/94	11/18/94	11/18/94

ANALYTE	SOW-3/90 - 11	CRQL				
3-Nitroaniline	25	25 U	25 U	25 U	25 U	25 U
Acenaphthene	10	10 U	10 U	10 U	10 U	10 U
2,4-Dinitrophenol	25	25 U	25 U	25 U	25 U	25 U
4-Nitrophenol	25	25 U	25 U	25 U	25 U	25 U
Dibenzofuran	10	10 U	10 U	10 U	10 U	10 U
2,4-Dinitrotoluene	10	10 U	10 U	10 U	10 U	10 U
Diethylphthalate	10	10 U	10 U	10 U	10 U	10 U
4-Chlorophenyl-phenylether	10	10 U	10 U	10 U	10 U	10 U
Fluorene	10	10 U	10 U	10 U	10 U	10 U
4-Nitroaniline	25	25 U	25 U	25 U	25 U	25 U
4,6-Dinitro-2-methylphenol	25	25 U	25 U	25 U	25 U	25 U
N-Nitrosodiphenylamine	10	10 U	10 U	10 U	10 U	10 U
4-Bromophenyl-phenylether	10	10 U	10 U	10 U	10 U	10 U
Hexachlorobenzene	10	10 U	10 U	10 U	10 U	10 U
Pentachlorophenol	25	25 U	25 U	25 U	25 U	25 U
Phenanthrene	10	10 U	10 U	10 U	10 U	10 U
Anthracene	10	10 U	10 U	10 U	10 U	10 U
Carbazole	10	10 U	10 U	10 U	10 U	10 U
Di-n-butylphthalate	10	10 U	10 U	10 U	10 U	10 U
Fluoranthene	10	10 U	10 U	10 U	10 U	10 U
Pyrene	10	10 U	10 U	10 U	10 U	10 U
Butylbenzylphthalate	10	10 U	10 U	10 U	10 U	10 U
3,3'-Dichlorobenzidine	10	10 U	10 U	10 U	10 U	10 U
Benzo(a)Anthracene	10	10 U	10 U	10 U	10 U	10 U
Chrysene	10	10 U	10 U	10 U	10 U	10 U
bis(2-Ethylhexyl)phthalate	10	2 J	10 U	3 JB	2 JB	2 JB
Di-n-octylphthalate	10	10 U	10 U	10 U	10 U	10 U
Benzo(b)Fluoranthene	10	10 U	10 U	10 U	10 U	10 U
Benzo(k)Fluoranthene	10	10 U	10 U	10 U	10 U	10 U
Benzo(a)Pyrene	10	10 U	10 U	10 U	10 U	10 U
Indeno(1,2,3-c,d)Pyrene	10	10 U	10 U	10 U	10 U	10 U
Dibenz(a,h)Anthracene	10	10 U	10 U	10 U	10 U	10 U
Benzo(g,h,i)perylene	10	10 U	10 U	10 U	10 U	10 U
=====						
Dilution Factor:	1.00	1.00	1.00	1.00	1.00	1.00
Sample Volume\Weight (ml\g):	1000	1000	1000	1000	1000	1000
Associated Method Blank:	R1695.D	S2097.D	S1523.D	S1523.D		
Associated Equipment Blank:	-	-	-	-	-	-
Associated Field Blank:	-	-	-	-	-	-

Site: EQUIPMENT RINSATE

U: not detected J: estimated B: blank contamination

Table 1
Laboratory Report of Analysis

ANALYTE	LOCATION: CL-102 DUP CL-102 CL-103 CL-103 CL-104 CL-104						
	ISIS ID: ENCL102XXX94XD ENCL102XXX94XX ENCL103XXX94XX ENCL103XXX94XX ENCL104XXX94XX ENCL104XXX94XX						
	LAB NUMBER: 2223317 2223314 2223318 2223318 R 2223312 2223312 D						
	DATE SAMPLED: 10/04/94 10/04/94 10/04/94 10/04/94 10/05/94 10/05/94						
	DATE EXTRACTED: 10/11/94 10/11/94 10/11/94 10/11/94 10/11/94 10/11/94						
	DATE ANALYZED: 11/18/94 11/18/94 11/18/94 11/18/94 11/18/94 11/21/94						
ANALYTE	SOW-3/90 - II	CRQL					
Phenol	10	50 U	50 U	50 U	50 U	50 U	200 U
bis(2-Chloroethyl)ether	10	50 U	50 U	50 U	50 U	50 U	200 U
2-Chlorophenol	10	50 U	50 U	50 U	50 U	50 U	200 U
1,3-Dichlorobenzene	10	50 U	50 U	50 U	50 U	50 U	200 U
1,4-Dichlorobenzene	10	50 U	50 U	50 U	50 U	50 U	200 U
1,2-Dichlorobenzene	10	50 U	50 U	50 U	50 U	9 J	200 U
2-Methylphenol	10	50 U	50 U	50 U	50 U	50 U	200 U
2,2'-oxybis(1-Chloropropane)	10	50 U	50 U	50 U	50 U	50 U	200 U
4-Methylphenol	10	50 U	50 U	50 U	50 U	32 J	31 JD
N-Nitroso-di-n-propylamine	10	50 U	50 U	50 U	50 U	50 U	200 U
Hexachloroethane	10	50 U	50 U	50 U	50 U	50 U	200 U
Nitrobenzene	10	50 U	50 U	50 U	50 U	50 U	200 U
Isophorone	10	50 U	50 U	50 U	50 U	50 U	200 U
2-Nitrophenol	10	50 U	50 U	50 U	50 U	50 U	200 U
2,4-Dimethylphenol	10	50 U	50 U	50 U	50 U	50 U	200 U
bis(2-Chloroethoxy)methane	10	50 U	50 U	50 U	50 U	50 U	200 U
2,4-Dichlorophenol	10	50 U	50 U	50 U	50 U	50 U	200 U
1,2,4-Trichlorobenzene	10	50 U	50 U	50 U	50 U	50 U	200 U
Naphthalene	10	50 U	50 U	50 U	50 U	14 J	200 U
4-Chloroaniline	10	50 U	50 U	50 U	50 U	50 U	200 U
Hexachlorobutadiene	10	50 U	50 U	50 U	50 U	50 U	200 U
4-Chloro-3-Methylphenol	10	50 U	50 U	50 U	50 U	50 U	200 U
2-Methylnaphthalene	10	50 U	50 U	50 U	50 U	28 J	22 JD
Hexachlorocyclopentadiene	10	50 U	50 U	50 U	50 U	50 U	200 U
2,4,6-Trichlorophenol	10	50 U	50 U	50 U	50 U	50 U	200 U
2,4,5-Trichlorophenol	25	120 U	120 U	120 U	120 U	120 U	500 U
2-Chloronaphthalene	10	50 U	50 U	50 U	50 U	50 U	200 U
2-Nitroaniline	25	120 U	120 U	120 U	120 U	120 U	500 U
Dimethylphthalate	10	50 U	50 U	50 U	5 J	50 U	200 U
Acenaphthylene	10	50 U	50 U	50 U	50 U	50 U	200 U
2,6-Dinitrotoluene	10	50 U	50 U	50 U	50 U	50 U	200 U

Site: SUMP LIQUIDS

U: not detected J: estimated D: diluted result B: blank contamination

Table 1
Laboratory Report of Analysis

LOCATION:	CL-102 DUP	CL-102	CL-103	CL-103	CL-104	CL-104
ISIS ID:	ENCL102XXX94XD	ENCL102XXX94XX	ENCL103XXX94XX	ENCL103XXX94XX	ENCL104XXX94XX	ENCL104XXX94XX
LAB NUMBER:	2223317	2223314	2223318	2223318 R	2223312	2223312 D
DATE SAMPLED:	10/04/94	10/04/94	10/04/94	10/04/94	10/05/94	10/05/94
DATE EXTRACTED:	10/11/94	10/11/94	10/11/94	10/11/94	10/11/94	10/11/94
DATE ANALYZED:	11/18/94	11/18/94	11/18/94	11/18/94	11/18/94	11/21/94

ANALYTE	SOW-3/90 - II	CRQL						
3-Nitroaniline	25	120 U	120 U	120 U	120 U	120 U	500 U	
Acenaphthene	10	50 U	50 U	50 U	50 U	50 U	200 U	
2,4-Dinitrophenol	25	120 U	120 U	120 U	120 U	120 U	500 U	
4-Nitrophenol	25	120 U	120 U	120 U	120 U	120 U	500 U	
Dibenzofuran	10	50 U	50 U	50 U	50 U	50 U	200 U	
2,4-Dinitrotoluene	10	50 U	50 U	50 U	50 U	50 U	200 U	
Diethylphthalate	10	50 U	50 U	50 U	7 J	50 U	200 U	
4-Chlorophenyl-phenylether	10	50 U	50 U	50 U	50 U	50 U	200 U	
Fluorene	10	50 U	50 U	50 U	50 U	50 U	200 U	
4-Nitroaniline	25	120 U	120 U	120 U	120 U	120 U	500 U	
4,6-Dinitro-2-methylphenol	25	120 U	120 U	120 U	120 U	120 U	500 U	
N-Nitrosodiphenylamine	10	50 U	50 U	50 U	50 U	50 U	200 U	
4-Bromophenyl-phenylether	10	50 U	50 U	50 U	50 U	50 U	200 U	
Hexachlorobenzene	10	50 U	50 U	50 U	50 U	50 U	200 U	
Pentachlorophenol	25	120 U	120 U	120 U	120 U	560 E	180 JD	
Phenanthrene	10	50 U	50 U	50 U	50 U	50 U	200 U	
Anthracene	10	50 U	50 U	50 U	50 U	50 U	200 U	
Carbazole	10	50 U	50 U	50 U	50 U	50 U	200 U	
Di-n-butylphthalate	10	50 U	50 U	50 U	50 U	50 U	200 U	
Fluoranthene	10	50 U	50 U	50 U	50 U	50 U	200 U	
Pyrene	10	6 J	50 U	50 U	50 U	50 U	200 U	
Butylbenzylphthalate	10	38 J	38 J	50 U	50 U	20 J	30 JD	
3,3'-Dichlorobenzidine	10	50 U	50 U	50 U	50 U	50 U	200 U	
Benzo(a)Anthracene	10	50 U	50 U	50 U	50 U	50 U	200 U	
Chrysene	10	50 U	50 U	50 U	50 U	50 U	200 U	
bis(2-Ethylhexyl)phthalate	10	380 B	290 B	11 JB	20 JB	330 B	400 BD	
Di-n-octylphthalate	10	130	40 J	50 U	8 J	15 J	63 JD	
Benzo(b)Fluoranthene	10	50 U	50 U	50 U	50 U	50 U	200 U	
Benzo(k)Fluoranthene	10	50 U	50 U	50 U	50 U	50 U	200 U	
Benzo(a)Pyrene	10	50 U	50 U	50 U	50 U	50 U	200 U	
Indeno(1,2,3-c,d)Pyrene	10	50 U	50 U	50 U	50 U	50 U	200 U	
Dibenz(a,h)Anthracene	10	50 U	50 U	50 U	50 U	50 U	200 U	
Benzo(g,h,i)perylene	10	50 U	50 U	50 U	50 U	50 U	200 U	

	Dilution Factor:	5.00	5.00	5.00	5.00	5.00	20.0
Sample Volume\Weight (ml\g):		1000	1000	1000	1000	1000	1000
Associated Method Blank:		S1523.D	S1523.D	S1523.D	S1523.D	S1523.D	S1523.D
Associated Equipment Blank:		-	-	-	-	-	-
Associated Field Blank:		-	-	-	-	-	-

Site: SUMP LIQUIDS

U: not detected J: estimated D: diluted result B: blank contamination

Table 1
Laboratory Report of Analysis

LOCATION:	CD-101	CD-103
ISIS ID:	ENCD101XXX94XX	ENCD103XXX94XX
LAB NUMBER:	2223303	2223304
DATE SAMPLED:	10/04/94	10/04/94
DATE EXTRACTED:	10/08/94	10/08/94
DATE ANALYZED:	11/17/94	11/17/94

ANALYTE	SOW-3/90 - 11	CRQL		
Phenol	330	21000	U	30000 U
bis(2-Chloroethyl)ether	330	21000	U	30000 U
2-Chlorophenol	330	21000	U	30000 U
1,3-Dichlorobenzene	330	21000	U	30000 U
1,4-Dichlorobenzene	330	21000	U	30000 U
1,2-Dichlorobenzene	330	21000	U	30000 U
2-Methylphenol	330	21000	U	30000 U
2,2'-oxybis(1-Chloropropane)	330	21000	U	30000 U
4-Methylphenol	330	21000	U	30000 U
N-Nitroso-di-n-propylamine	330	21000	U	30000 U
Hexachloroethane	330	21000	U	30000 U
Nitrobenzene	330	21000	U	30000 U
Isophorone	330	21000	U	30000 U
2-Nitrophenol	330	21000	U	30000 U
2,4-Dimethylphenol	330	21000	U	30000 U
bis(2-Chloroethoxy)methane	330	21000	U	30000 U
2,4-Dichlorophenol	330	21000	U	30000 U
1,2,4-Trichlorobenzene	330	21000	U	30000 U
Naphthalene	330	21000	U	30000 U
4-Chloroaniline	330	21000	U	30000 U
Hexachlorobutadiene	330	21000	U	30000 U
4-Chloro-3-Methylphenol	330	21000	U	30000 U
2-Methylnaphthalene	330	21000	U	30000 U
Hexachlorocyclopentadiene	330	21000	U	30000 U
2,4,6-Trichlorophenol	330	21000	U	30000 U
2,4,5-Trichlorophenol	800	50000	U	73000 U
2-Chloronaphthalene	330	21000	U	30000 U
2-Nitroaniline	800	50000	U	73000 U
Dimethylphthalate	330	21000	U	30000 U
Acenaphthylene	330	21000	U	30000 U
2,6-Dinitrotoluene	330	21000	U	30000 U

Site: SUMP SEDIMENTS

U: not detected J: estimated

Table 1
Laboratory Report of Analysis

LOCATION:	CD-101	CD-103
ISIS ID:	ENCD101XXX94XX	ENCD103XXX94XX
LAB NUMBER:	2223303	2223304
DATE SAMPLED:	10/04/94	10/04/94
DATE EXTRACTED:	10/08/94	10/08/94
DATE ANALYZED:	11/17/94	11/17/94

ANALYTE	SOW-3/90 - 11	CRQL				
3-Nitroaniline	800	50000	U	73000	U	
Acenaphthene	330	21000	U	30000	U	
2,4-Dinitrophenol	800	50000	U	73000	U	
4-Nitrophenol	800	50000	U	73000	U	
Dibenzofuran	330	21000	U	30000	U	
2,4-Dinitrotoluene	330	21000	U	30000	U	
Diethylphthalate	330	21000	U	30000	U	
4-Chlorophenyl-phenylether	330	21000	U	30000	U	
Fluorene	330	21000	U	30000	U	
4-Nitroaniline	800	50000	U	73000	U	
4,6-Dinitro-2-methylphenol	800	50000	U	73000	U	
N-Nitrosodiphenylamine	330	21000	U	30000	U	
4-Bromophenyl-phenylether	330	21000	U	30000	U	
Hexachlorobenzene	330	21000	U	30000	U	
Pentachlorophenol	800	50000	U	73000	U	
Phenanthrene	330	21000	U	30000	U	
Anthracene	330	21000	U	30000	U	
Carbazole	330	21000	U	30000	U	
Di-n-butylphthalate	330	21000	U	30000	U	
Fluoranthene	330	21000	U	30000	U	
Pyrene	330	21000	U	30000	U	
Butylbenzylphthalate	330	3000	J	30000	U	
3,3'-Dichlorobenzidine	330	21000	U	30000	U	
Benzo(a)Anthracene	330	21000	U	30000	U	
Chrysene	330	21000	U	30000	U	
bis(2-Ethylhexyl)phthalate	330	46000		6900	J	
Di-n-octylphthalate	330	21000	U	30000	U	
Benzo(b)Fluoranthene	330	21000	U	30000	U	
Benzo(k)Fluoranthene	330	21000	U	30000	U	
Benzo(a)Pyrene	330	21000	U	30000	U	
Indeno(1,2,3-c,d)Pyrene	330	21000	U	30000	U	
Dibenz(a,h)Anthracene	330	21000	U	30000	U	
Benzo(g,h,i)perylene	330	21000	U	30000	U	

Dilution Factor:	20.0	50.0
Percent Solids:	32	55
Sample Volume\Weight (ml\g):	30.0	30.0

Associated Method Blank:	S1503.D	S1503.D
Associated Equipment Blank:	ENQXXX1XXX94XX	ENQXXX1XXX94XX
Associated Field Blank:	-	-

Site: SUMP SEDIMENTS

U: not detected J: estimated

Table 1
Laboratory Report of Analysis

ANALYTE	LOCATION: WT-101 DUP WT-101 WT-102 WT-103 WT-104 WT-105 WT-106 WT-106								
	ISIS ID: ENWT101XXX94XD ENWT101XXX94XX ENWT102XXX94XX ENWT103XXX94XX ENWT104XXX94XX ENWT105XXX94XX ENWT106XXX94XX ENWT106XXX94XX								
	LAB NUMBER: 2223309 2223306 2223305 2223310 2223311 2223301 2223302 2223302 R								
	DATE SAMPLED: 10/04/94 10/04/94 10/04/94 10/04/94 10/04/94 10/05/94 10/05/94 10/05/94								
	DATE EXTRACTED: 10/08/94 10/08/94 10/08/94 10/08/94 10/08/94 10/08/94 10/08/94 10/08/94								
	DATE ANALYZED: 11/17/94 11/17/94 11/17/94 11/17/94 11/17/94 11/17/94 11/17/94 11/17/94								
ANALYTE	SOW-3/90 - 11	CRQL							
Phenol	330	7900 U	150000 U	69000 U	11000 U	11000 U	7300 U	37000 U	37000 U
bis(2-Chloroethyl)ether	330	7900 U	150000 U	69000 U	11000 U	11000 U	7300 U	37000 U	37000 U
2-Chlorophenol	330	7900 U	150000 U	69000 U	11000 U	11000 U	7300 U	37000 U	37000 U
1,3-Dichlorobenzene	330	7900 U	150000 U	69000 U	11000 U	11000 U	7300 U	37000 U	37000 U
1,4-Dichlorobenzene	330	7900 U	150000 U	69000 U	11000 U	11000 U	7300 U	37000 U	37000 U
1,2-Dichlorobenzene	330	7900 U	150000 U	69000 U	11000 U	11000 U	7300 U	37000 U	37000 U
2-Methylphenol	330	7900 U	150000 U	69000 U	11000 U	11000 U	7300 U	37000 U	37000 U
2,2'-oxybis(1-Chloropropane)	330	7900 U	150000 U	69000 U	11000 U	11000 U	7300 U	37000 U	37000 U
4-Methylphenol	330	7900 U	150000 U	69000 U	11000 U	11000 U	7300 U	37000 U	37000 U
N-Nitroso-di-n-propylamine	330	7900 U	150000 U	69000 U	11000 U	11000 U	7300 U	37000 U	37000 U
Hexachloroethane	330	7900 U	150000 U	69000 U	11000 U	11000 U	7300 U	37000 U	37000 U
Nitrobenzene	330	7900 U	150000 U	69000 U	11000 U	11000 U	7300 U	37000 U	37000 U
Isophorone	330	7900 U	150000 U	69000 U	11000 U	11000 U	7300 U	37000 U	37000 U
2-Nitrophenol	330	7900 U	150000 U	69000 U	11000 U	11000 U	7300 U	37000 U	37000 U
2,4-Dimethylphenol	330	7900 U	150000 U	69000 U	11000 U	11000 U	7300 U	37000 U	37000 U
bis(2-Chloroethoxy)methane	330	7900 U	150000 U	69000 U	11000 U	11000 U	7300 U	37000 U	37000 U
2,4-Dichlorophenol	330	7900 U	150000 U	69000 U	11000 U	11000 U	7300 U	37000 U	37000 U
1,2,4-Trichlorobenzene	330	7900 U	150000 U	69000 U	11000 U	11000 U	7300 U	37000 U	37000 U
Naphthalene	330	7900 U	150000 U	69000 U	11000 U	11000 U	7300 U	37000 U	37000 U
4-Chloroaniline	330	7900 U	150000 U	69000 U	11000 U	11000 U	7300 U	37000 U	37000 U
Hexachlorobutadiene	330	7900 U	150000 U	69000 U	11000 U	11000 U	7300 U	37000 U	37000 U
4-Chloro-3-Methylphenol	330	7900 U	150000 U	69000 U	11000 U	11000 U	7300 U	37000 U	37000 U
2-Methylnaphthalene	330	7900 U	150000 U	69000 U	11000 U	11000 U	7300 U	37000 U	37000 U
Hexachlorocyclopentadiene	330	7900 U	150000 U	69000 U	11000 U	11000 U	7300 U	37000 U	37000 U
2,4,6-Trichlorophenol	330	7900 U	150000 U	69000 U	11000 U	11000 U	7300 U	37000 U	37000 U
2,4,5-Trichlorophenol	800	19000 U	360000 U	170000 U	27000 U	27000 U	18000 U	88000 U	88000 U
2-Chloronaphthalene	330	7900 U	150000 U	69000 U	11000 U	11000 U	7300 U	37000 U	37000 U
2-Nitroaniline	800	19000 U	360000 U	170000 U	27000 U	27000 U	18000 U	88000 U	88000 U
Dimethylphthalate	330	7900 U	150000 U	69000 U	11000 U	11000 U	7300 U	37000 U	37000 U
Acenaphthylene	330	7900 U	150000 U	69000 U	11000 U	11000 U	7300 U	37000 U	37000 U
2,6-Dinitrotoluene	330	7900 U	150000 U	69000 U	11000 U	11000 U	7300 U	37000 U	37000 U

Site: WASTE-SLUDGE

U: not detected J: estimated

Table 1
Laboratory Report of Analysis

	LOCATION:	WT-101 DUP	WT-101	WT-102	WT-103	WT-104	WT-105	WT-106	WT-106
	ISIS ID:	ENWT101XXX94XD	ENWT101XXX94XX	ENWT102XXX94XX	ENWT103XXX94XX	ENWT104XXX94XX	ENWT105XXX94XX	ENWT106XXX94XX	ENWT106XXX94XX
	LAB NUMBER:	2223309	2223306	2223305	2223310	2223311	2223301	2223302	2223302 R
	DATE SAMPLED:	10/04/94	10/04/94	10/04/94	10/04/94	10/04/94	10/05/94	10/05/94	10/05/94
	DATE EXTRACTED:	10/08/94	10/08/94	10/08/94	10/08/94	10/08/94	10/08/94	10/08/94	10/08/94
	DATE ANALYZED:	11/17/94	11/17/94	11/17/94	11/17/94	11/17/94	11/17/94	11/17/94	11/17/94
ANALYTE	SOW-3/90 - II	CRQL							
3-Nitroaniline	800	19000 U	360000 U	170000 U	27000 U	27000 U	18000 U	88000 U	88000 U
Acenaphthene	330	7900 U	150000 U	69000 U	11000 U	11000 U	7300 U	37000 U	37000 U
2,4-Dinitrophenol	800	19000 U	360000 U	170000 U	27000 U	27000 U	18000 U	88000 U	88000 U
4-Nitrophenol	800	19000 U	360000 U	170000 U	27000 U	27000 U	18000 U	88000 U	88000 U
Dibenzofuran	330	7900 U	150000 U	69000 U	11000 U	11000 U	7300 U	37000 U	37000 U
2,4-Dinitrotoluene	330	7900 U	150000 U	69000 U	11000 U	11000 U	7300 U	37000 U	37000 U
Diethylphthalate	330	7900 U	150000 U	69000 U	11000 U	11000 U	7300 U	37000 U	37000 U
4-Chlorophenyl-phenylether	330	7900 U	150000 U	69000 U	11000 U	11000 U	7300 U	37000 U	37000 U
Fluorene	330	7900 U	150000 U	69000 U	11000 U	11000 U	7300 U	37000 U	37000 U
4-Nitroaniline	800	19000 U	360000 U	170000 U	27000 U	27000 U	18000 U	88000 U	88000 U
4,6-Dinitro-2-methylphenol	800	19000 U	360000 U	170000 U	27000 U	27000 U	18000 U	88000 U	88000 U
N-Nitrosodiphenylamine	330	7900 U	150000 U	69000 U	11000 U	11000 U	7300 U	37000 U	37000 U
4-Bromophenyl-phenylether	330	7900 U	150000 U	69000 U	11000 U	11000 U	7300 U	37000 U	37000 U
Hexachlorobenzene	330	7900 U	150000 U	69000 U	11000 U	11000 U	7300 U	37000 U	37000 U
Pentachlorophenol	800	19000 U	360000 U	170000 U	27000 U	27000 U	18000 U	7400 J	8300 J
Phenanthrene	330	1100 J	150000 U	69000 U	11000 U	11000 U	7300 U	37000 U	37000 U
Anthracene	330	7900 U	150000 U	69000 U	11000 U	11000 U	7300 U	37000 U	37000 U
Carbazole	330	7900 U	150000 U	69000 U	11000 U	11000 U	7300 U	37000 U	37000 U
Di-n-butylphthalate	330	7900 U	150000 U	69000 U	11000 U	11000 U	7300 U	37000 U	37000 U
Fluoranthene	330	940 J	150000 U	69000 U	11000 U	11000 U	7300 U	37000 U	37000 U
Pyrene	330	1800 J	150000 U	69000 U	11000 U	11000 U	7300 U	37000 U	37000 U
Butylbenzylphthalate	330	7900 U	150000 U	69000 U	11000 U	11000 U	7300 U	37000 U	37000 U
3,3'-Dichlorobenzidine	330	7900 U	150000 U	69000 U	11000 U	11000 U	7300 U	37000 U	37000 U
Benzo(a)Anthracene	330	7900 U	150000 U	69000 U	11000 U	11000 U	7300 U	37000 U	37000 U
Chrysene	330	940 J	150000 U	69000 U	11000 U	11000 U	7300 U	37000 U	37000 U
bis(2-Ethylhexyl)phthalate	330	4400 J	25000 J	7900 J	11000 U	1900 J	7300 U	37000 U	37000 U
Di-n-octylphthalate	330	7900 U	150000 U	69000 U	11000 U	11000 U	7300 U	37000 U	37000 U
Benzo(b)Fluoranthene	330	7900 U	150000 U	69000 U	11000 U	11000 U	7300 U	37000 U	37000 U
Benzo(k)Fluoranthene	330	7900 U	150000 U	69000 U	11000 U	11000 U	7300 U	37000 U	37000 U
Benzo(a)Pyrene	330	7900 U	150000 U	69000 U	11000 U	11000 U	7300 U	37000 U	37000 U
Indeno(1,2,3-c,d)Pyrene	330	7900 U	150000 U	69000 U	11000 U	11000 U	7300 U	37000 U	37000 U
Dibenz(a,h)Anthracene	330	7900 U	150000 U	69000 U	11000 U	11000 U	7300 U	37000 U	37000 U
Benzo(g,h,i)perylene	330	7900 U	150000 U	69000 U	11000 U	11000 U	7300 U	37000 U	37000 U
=====									
Dilution Factor:	20.0	400	100	20.0	20.0	20.0	100	100	
Percent Solids:	84	88	48	59	60	91	91	91	
Sample Volume\Weight (ml\g):	30.0	30.0	30.0	30.0	30.0	30.0	30.0	30.0	
Associated Method Blank:	S1503.D	S1503.D	S1503.D	S1503.D	S1503.D	S1503.D	S1503.D	S1503.D	S1503.D
Associated Equipment Blank:	ENQSXX1XXX94XX	ENQSXX1XXX94XX	ENQSXX1XXX94XX	ENQSXX1XXX94XX	ENQSXX1XXX94XX	ENQSXX1XXX94XX	ENQSXX1XXX94XX	ENQSXX1XXX94XX	ENQSXX1XXX94XX
Associated Field Blank:	-	-	-	-	-	-	-	-	-

Site: WASTE-SLUDGE

U: not detected J: estimated

Table 1
Laboratory Report of Analysis

LOCATION:	SS-101	SS-102 DUP	SS-102
ISIS ID:	ENSS101XXX94XX	ENSS102XXX94XD	ENSS102XXX94XX
LAB NUMBER:	2241101	2241105	2241102
DATE SAMPLED:	10/28/94	10/28/94	10/28/94
DATE EXTRACTED:	11/01/94	11/01/94	11/01/94
DATE ANALYZED:	12/03/94	12/03/94	12/03/94

ANALYTE	SOW-3/90 - II	CRQL			
Phenol	330	1800 U	900 U	900 U	
bis(2-Chloroethyl)ether	330	1800 U	900 U	900 U	
2-Chlorophenol	330	1800 U	900 U	900 U	
1,3-Dichlorobenzene	330	1800 U	900 U	900 U	
1,4-Dichlorobenzene	330	1800 U	900 U	900 U	
1,2-Dichlorobenzene	330	1800 U	900 U	900 U	
2-Methylphenol	330	1800 U	900 U	900 U	
2,2'-oxybis(1-Chloropropane)	330	1800 U	900 U	900 U	
4-Methylphenol	330	1800 U	900 U	900 U	
N-Nitroso-di-n-propylamine	330	1800 U	900 U	900 U	
Hexachloroethane	330	1800 U	900 U	900 U	
Nitrobenzene	330	1800 U	900 U	900 U	
Isophorone	330	1800 U	900 U	900 U	
2-Nitrophenol	330	1800 U	900 U	900 U	
2,4-Dimethylphenol	330	1800 U	900 U	900 U	
bis(2-Chloroethoxy)methane	330	1800 U	900 U	900 U	
2,4-Dichlorophenol	330	1800 U	900 U	900 U	
1,2,4-Trichlorobenzene	330	1800 U	900 U	900 U	
Naphthalene	330	1800 U	900 U	900 U	
4-Chloroaniline	330	1800 U	900 U	900 U	
Hexachlorobutadiene	330	1800 U	900 U	900 U	
4-Chloro-3-Methylphenol	330	1800 U	900 U	900 U	
2-Methylnaphthalene	330	1800 U	900 U	900 U	
Hexachlorocyclopentadiene	330	1800 U	900 U	900 U	
2,4,6-Trichlorophenol	330	1800 U	900 U	900 U	
2,4,5-Trichlorophenol	800	4400 U	2200 U	2200 U	
2-Chloronaphthalene	330	1800 U	900 U	900 U	
2-Nitroaniline	800	4400 U	2200 U	2200 U	
Dimethylphthalate	330	1800 U	900 U	900 U	
Acenaphthylene	330	1800 U	900 U	900 U	
2,6-Dinitrotoluene	330	1800 U	900 U	900 U	

Site: SURFACE SOILS

U: not detected J: estimated

Table 1
Laboratory Report of Analysis

LOCATION:	SS-101	SS-102 DUP	SS-102
ISIS ID:	ENSS101XXX94XX	ENSS102XXX94XD	ENSS102XXX94XX
LAB NUMBER:	2241101	2241105	2241102
DATE SAMPLED:	10/28/94	10/28/94	10/28/94
DATE EXTRACTED:	11/01/94	11/01/94	11/01/94
DATE ANALYZED:	12/03/94	12/03/94	12/03/94

ANALYTE	SOW-3/90 - 11	CRQL			
3-Nitroaniline	800	4400 U	2200 U	2200 U	
Acenaphthene	330	470 J	900 U	900 U	
2,4-Dinitrophenol	800	4400 U	2200 U	2200 U	
4-Nitrophenol	800	4400 U	2200 U	2200 U	
Dibenzofuran	330	290 J	900 U	900 U	
2,4-Dinitrotoluene	330	1800 U	900 U	900 U	
Diethylphthalate	330	1800 U	900 U	900 U	
4-Chlorophenyl-phenylether	330	1800 U	900 U	900 U	
Fluorene	330	560 J	900 U	900 U	
4-Nitroaniline	800	4400 U	2200 U	2200 U	
4,6-Dinitro-2-methylphenol	800	4400 U	2200 U	2200 U	
N-Nitrosodiphenylamine	330	1800 U	900 U	900 U	
4-Bromophenyl-phenylether	330	1800 U	900 U	900 U	
Hexachlorobenzene	330	1800 U	900 U	900 U	
Pentachlorophenol	800	4400 U	2200 U	2200 U	
Phenanthrene	330	6100	220 J	450 J	
Anthracene	330	1100 J	900 U	900 U	
Carbazole	330	640 J	900 U	900 U	
Di-n-butylphthalate	330	1800 U	900 U	900 U	
Fluoranthene	330	9400	320 J	680 J	
Pyrene	330	7900	300 J	620 J	
Butylbenzylphthalate	330	1800 U	900 U	900 U	
3,3'-Dichlorobenzidine	330	1800 U	900 U	900 U	
Benzo(a)Anthracene	330	3900	160 J	350 J	
Chrysene	330	4400	200 J	450 J	
bis(2-Ethylhexyl)phthalate	330	1800 U	900 U	900 U	
Di-n-octylphthalate	330	1800 U	900 U	900 U	
Benzo(b)Fluoranthene	330	3200	130 J	270 J	
Benzo(k)Fluoranthene	330	2100	900 U	170 J	
Benzo(a)Pyrene	330	2400	95 J	210 J	
Indeno(1,2,3-c,d)Pyrene	330	720 J	900 U	120 J	
Dibenz(a,h)Anthracene	330	1800 U	900 U	900 U	
Benzo(g,h,i)perylene	330	600 J	900 U	120 J	

Dilution Factor:	5.00	2.00	2.00
Percent Solids:	90	74	74
Sample Volume\Weight (ml\g):	30.0	30.0	30.0
Associated Method Blank:	R1715.D	R1715.D	R1715.D
Associated Equipment Blank:	ENQS003XXX94XX	ENQS003XXX94XX	ENQS003XXX94XX
Associated Field Blank:			

Site: SURFACE SOILS

U: not detected J: estimated

Table 1
Laboratory Report of Analysis

LOCATION:	BS-101	BS-102
ISIS ID:	ENBS101XX694XX	ENBS102XX694XX
LAB NUMBER:	2241106	2241107
DATE SAMPLED:	10/27/94	10/28/94
DATE EXTRACTED:	11/01/94	11/01/94
DATE ANALYZED:	12/03/94	12/03/94

ANALYTE	SOW-3/90 - II	CRQL		
Phenol	330	860	U	410 U
bis(2-Chloroethyl)ether	330	860	U	410 U
2-Chlorophenol	330	860	U	410 U
1,3-Dichlorobenzene	330	860	U	410 U
1,4-Dichlorobenzene	330	860	U	410 U
1,2-Dichlorobenzene	330	860	U	410 U
2-Methylphenol	330	860	U	410 U
2,2'-oxybis(1-Chloropropane)	330	860	U	410 U
4-Methylphenol	330	860	U	410 U
N-Nitroso-di-n-propylamine	330	860	U	410 U
Hexachloroethane	330	860	U	410 U
Nitrobenzene	330	860	U	410 U
Isophorone	330	860	U	410 U
2-Nitrophenol	330	860	U	410 U
2,4-Dimethylphenol	330	860	U	410 U
bis(2-Chloroethoxy)methane	330	860	U	410 U
2,4-Dichlorophenol	330	860	U	410 U
1,2,4-Trichlorobenzene	330	860	U	410 U
Naphthalene	330	860	U	410 U
4-Chloroaniline	330	860	U	410 U
Hexachlorobutadiene	330	860	U	410 U
4-Chloro-3-Methylphenol	330	860	U	410 U
2-Methylnaphthalene	330	860	U	410 U
Hexachlorocyclopentadiene	330	860	U	410 U
2,4,6-Trichlorophenol	330	860	U	410 U
2,4,5-Trichlorophenol	800	2100	U	990 U
2-Chloronaphthalene	330	860	U	410 U
2-Nitroaniline	800	2100	U	990 U
Dimethylphthalate	330	860	U	410 U
Acenaphthylene	330	860	U	410 U
2,6-Dinitrotoluene	330	860	U	410 U

Site: SOIL BORINGS

U: not detected J: estimated

Table 1
Laboratory Report of Analysis

LOCATION:	BS-101	BS-102
ISIS ID:	ENBS101XX694XX	ENBS102XX694XX
LAB NUMBER:	2241106	2241107
DATE SAMPLED:	10/27/94	10/28/94
DATE EXTRACTED:	11/01/94	11/01/94
DATE ANALYZED:	12/03/94	12/03/94

ANALYTE	SOW-3/90 - 11	CRQL		
3-Nitroaniline	800	2100	U	990 U
Acenaphthene	330	860	U	410 U
2,4-Dinitrophenol	800	2100	U	990 U
4-Nitrophenol	800	2100	U	990 U
Dibenzofuran	330	860	U	410 U
2,4-Dinitrotoluene	330	860	U	410 U
Diethylphthalate	330	860	U	410 U
4-Chlorophenyl-phenylether	330	860	U	410 U
Fluorene	330	860	U	410 U
4-Nitroaniline	800	2100	U	990 U
4,6-Dinitro-2-methylphenol	800	2100	U	990 U
N-Nitrosodiphenylamine	330	860	U	410 U
4-Bromophenyl-phenylether	330	860	U	410 U
Hexachlorobenzene	330	860	U	410 U
Pentachlorophenol	800	2100	U	990 U
Phenanthrene	330	300	J	110 J
Anthracene	330	860	U	410 U
Carbazole	330	860	U	410 U
Di-n-butylphthalate	330	860	U	410 U
Fluoranthene	330	260	J	200 J
Pyrene	330	240	J	180 J
Butylbenzylphthalate	330	860	U	410 U
3,3'-Dichlorobenzidine	330	860	U	410 U
Benzo(a)Anthracene	330	130	J	100 J
Chrysene	330	140	J	120 J
bis(2-Ethylhexyl)phthalate	330	860	U	110 J
Di-n-octylphthalate	330	860	U	410 U
Benzo(b)Fluoranthene	330	860	U	85 J
Benzo(k)Fluoranthene	330	860	U	61 J
Benzo(a)Pyrene	330	860	U	65 J
Indeno(1,2,3-c,d)Pyrene	330	860	U	44 J
Dibenz(a,h)Anthracene	330	860	U	410 U
Benzo(g,h,i)perylene	330	860	U	45 J

Dilution Factor:	2.00	1.00
Percent Solids:	77	81
Sample Volume\Weight (ml\g):	30.0	30.0

Associated Method Blank:	R1715.D	R1715.D
Associated Equipment Blank:	ENQS003XXX94XX	ENQS003XXX94XX
Associated Field Blank:	-	-

Site: SOIL BORINGS
U: not detected J: estimated

Table 1
Laboratory Report of Analysis

LOCATION:	MW-101 DUP	MW-101	MW-102	MW-103
ISIS ID:	ENMW101XXX94XD	ENMW101XXX94XX	ENMW102XXX94XX	ENMW103XXX94XX
LAB NUMBER:	2263804	2263801	2263805	2263806
DATE SAMPLED:	11/30/94	11/30/94	11/30/94	11/30/94
DATE EXTRACTED:	12/06/94	12/06/94	12/06/94	12/06/94
DATE ANALYZED:	12/22/94	12/22/94	12/22/94	12/22/94

ANALYTE	SOW-3/90 - II	CRQL				
Phenol	10		10 U	10 U	10 U	10 U
bis(2-Chloroethyl)ether	10		10 U	10 U	10 U	10 U
2-Chlorophenol	10		10 U	10 U	10 U	10 U
1,3-Dichlorobenzene	10		10 U	10 U	10 U	10 U
1,4-Dichlorobenzene	10		10 U	10 U	10 U	10 U
1,2-Dichlorobenzene	10		10 U	10 U	10 U	10 U
2-Methylphenol	10		10 U	10 U	10 U	2 J
2,2'-oxybis(1-Chloropropane)	10		10 U	10 U	10 U	10 U
4-Methylphenol	10		10 U	10 U	10 U	15
N-Nitroso-di-n-propylamine	10		10 U	10 U	10 U	10 U
Hexachloroethane	10		10 U	10 U	10 U	10 U
Nitrobenzene	10		10 U	10 U	10 U	10 U
Isophorone	10		10 U	10 U	10 U	10 U
2-Nitrophenol	10		10 U	10 U	10 U	10 U
2,4-Dimethylphenol	10		10 U	10 U	10 U	1 J
bis(2-Chloroethoxy)methane	10		10 U	10 U	10 U	10 U
2,4-Dichlorophenol	10		10 U	10 U	10 U	10 U
1,2,4-Trichlorobenzene	10		10 U	10 U	10 U	10 U
Naphthalene	10		10 U	10 U	10 U	14
4-Chloroaniline	10		10 U	10 U	10 U	10 U
Hexachlorobutadiene	10		10 U	10 U	10 U	10 U
4-Chloro-3-Methylphenol	10		10 U	10 U	10 U	10 U
2-Methylnaphthalene	10		10 U	10 U	10 U	4 J
Hexachlorocyclopentadiene	10		10 U	10 U	10 U	10 U
2,4,6-Trichlorophenol	10		10 U	10 U	10 U	10 U
2,4,5-Trichlorophenol	25		25 U	25 U	25 U	25 U
2-Chloronaphthalene	10		10 U	10 U	10 U	10 U
2-Nitroaniline	25		25 U	25 U	25 U	25 U
Dimethylphthalate	10		10 U	10 U	10 U	10 U
Acenaphthylene	10		10 U	10 U	10 U	10 U
2,6-Dinitrotoluene	10		10 U	10 U	10 U	10 U

Site: MONITORING WELLS

U: not detected J: estimated

Table 1
Laboratory Report of Analysis

LOCATION:	MW-101 DUP	MW-101	MW-102	MW-103
ISIS ID:	ENMW101XXX94XD	ENMW101XXX94XX	ENMW102XXX94XX	ENMW103XXX94XX
LAB NUMBER:	2263804	2263801	2263805	2263806
DATE SAMPLED:	11/30/94	11/30/94	11/30/94	11/30/94
DATE EXTRACTED:	12/06/94	12/06/94	12/06/94	12/06/94
DATE ANALYZED:	12/22/94	12/22/94	12/22/94	12/22/94

ANALYTE	SOW-3/90 - II	CRQL				
3-Nitroaniline	25	25 U	25 U	25 U	25 U	25 U
Acenaphthene	10	10 U	10 U	10 U	10 U	10 U
2,4-Dinitrophenol	25	25 U	25 U	25 U	25 U	25 U
4-Nitrophenol	25	25 U	25 U	25 U	25 U	25 U
Dibenzofuran	10	10 U	10 U	10 U	10 U	1 J
2,4-Dinitrotoluene	10	10 U	10 U	10 U	10 U	10 U
Diethylphthalate	10	10 U	10 U	10 U	10 U	10 U
4-Chlorophenyl-phenylether	10	10 U	10 U	10 U	10 U	10 U
Fluorene	10	10 U	10 U	10 U	10 U	2 J
4-Nitroaniline	25	25 U	25 U	25 U	25 U	25 U
4,6-Dinitro-2-methylphenol	25	25 U	25 U	25 U	25 U	25 U
N-Nitrosodiphenylamine	10	10 U	10 U	10 U	10 U	29
4-Bromophenyl-phenylether	10	10 U	10 U	10 U	10 U	10 U
Hexachlorobenzene	10	10 U	10 U	10 U	10 U	10 U
Pentachlorophenol	25	25 U	25 U	25 U	25 U	25 U
Phenanthrene	10	10 U	10 U	10 U	10 U	2 J
Anthracene	10	10 U	10 U	10 U	10 U	10 U
Carbazole	10	10 U	10 U	10 U	10 U	2 J
Di-n-butylphthalate	10	10 U	10 U	10 U	10 U	10 U
Fluoranthene	10	10 U	10 U	10 U	10 U	10 U
Pyrene	10	10 U	10 U	10 U	10 U	10 U
Butylbenzylphthalate	10	10 U	10 U	10 U	10 U	10 U
3,3'-Dichlorobenzidine	10	10 U	10 U	10 U	10 U	10 U
Benzo(a)Anthracene	10	10 U	10 U	10 U	10 U	10 U
Chrysene	10	10 U	10 U	10 U	10 U	10 U
bis(2-Ethylhexyl)phthalate	10	10 U	10 U	2 J	10 U	10 U
Di-n-octylphthalate	10	10 U	10 U	10 U	10 U	10 U
Benzo(b)Fluoranthene	10	10 U	10 U	10 U	10 U	10 U
Benzo(k)Fluoranthene	10	10 U	10 U	10 U	10 U	10 U
Benzo(a)Pyrene	10	10 U	10 U	10 U	10 U	10 U
Indeno(1,2,3-c,d)Pyrene	10	10 U	10 U	10 U	10 U	10 U
Dibenz(a,h)Anthracene	10	10 U	10 U	10 U	10 U	10 U
Benzo(g,h,i)perylene	10	10 U	10 U	10 U	10 U	10 U

Dilution Factor:	1.00	1.00	1.00	1.00
Sample Volume\Weight (ml\g):	1000	1000	1000	1000
Associated Method Blank:	S2097.D	S2097.D	S2097.D	S2097.D
Associated Equipment Blank:	ENQS004XXX94XX	ENQS004XXX94XX	ENQS004XXX94XX	ENQS004XXX94XX
Associated Field Blank:	-	-	-	-

Site: MONITORING WELLS
U: not detected J: estimated

Table 1
Laboratory Report of Analysis

LOCATION:	QS-003	QS-004	QS-XX1	QS-2
ISIS ID:	ENQS003XXX94XX	ENQS004XXX94XX	ENQSXX1XXX94XX	ENQSXX2XXX94XX
LAB NUMBER:	2241108	2263807	2223319	2223313
DATE SAMPLED:	10/28/94	11/30/94	10/04/94	10/04/94
DATE EXTRACTED:	11/01/94	12/03/94	10/08/94	10/08/94
DATE ANALYZED:	12/07/94	12/12/94	10/29/94	10/29/94

ANALYTE	SOW-3/90 - II	CRQL				
alpha-BHC	0.05	0.05 U	0.05 U	0.056 U	0.05 U	
beta-BHC	0.05	0.05 U	0.05 U	0.056 U	0.05 U	
delta-BHC	0.05	0.05 U	0.05 U	0.056 U	0.05 U	
gamma-BHC (Lindane)	0.05	0.05 U	0.05 U	0.056 U	0.05 U	
Heptachlor	0.05	0.05 U	0.05 U	0.056 U	0.05 U	
Aldrin	0.05	0.05 U	0.05 U	0.056 U	0.05 U	
Heptachlor Epoxide	0.05	0.05 U	0.05 U	0.056 U	0.05 U	
Endosulfan I	0.05	0.05 U	0.05 U	0.056 U	0.05 U	
Dieldrin	0.1	0.1 U	0.1 U	0.11 U	0.1 U	
4,4'-DDE	0.1	0.1 U	0.1 U	0.11 U	0.1 U	
Endrin	0.1	0.1 U	0.1 U	0.11 U	0.1 U	
Endosulfan II	0.1	0.1 U	0.1 U	0.11 U	0.1 U	
4,4'-DDD	0.1	0.1 U	0.1 U	0.11 U	0.1 U	
Endrin Aldehyde	0.1	0.1 U	0.1 U	0.11 U	0.1 U	
Endosulfan Sulfate	0.1	0.1 U	0.1 U	0.11 U	0.1 U	
4,4'-DDT	0.1	0.1 U	0.1 U	0.11 U	0.1 U	
Methoxychlor	0.5	0.5 U	0.5 U	0.56 U	0.5 U	
Endrin Ketone	0.1	0.1 U	0.1 U	0.11 U	0.1 U	
alpha-Chlordane	0.05	0.05 U	0.05 U	0.056 U	0.05 U	
gamma-Chlordane	0.05	0.05 U	0.05 U	0.056 U	0.05 U	
Toxaphene	5	5.0 U	5.0 U	5.6 U	5.0 U	
Aroclor-1016	1	1.0 U	1.0 U	1.1 U	1.0 U	
Aroclor-1221	2	2.0 U	2.0 U	2.2 U	2.0 U	
Aroclor-1232	1	1.0 U	1.0 U	1.1 U	1.0 U	
Aroclor-1242	1	1.0 U	1.0 U	1.1 U	1.0 U	
Aroclor-1248	1	1.0 U	1.0 U	1.1 U	1.0 U	
Aroclor-1254	1	1.0 U	1.0 U	1.1 U	1.0 U	
Aroclor-1260	1	1.0 U	1.0 U	1.1 U	1.0 U	

Dilution Factor:	1.00	1.00	1.00	1.00
Sample Volume/Weight (ml/g):	1000	1000	900	1000

Associated Method Blank:	PWB1101A1	PWB1203A	PWB1008B	PWB1008B
Associated Equipment Blank:	-	-	-	-
Associated Field Blank:	-	-	-	-

Site: EQUIPMENT RINSATE
U: not detected

Table 1
Laboratory Report of Analysis

	CL-102 DUP	CL-102	CL-103	CL-104	CL-104
LOCATION:	CL-102 DUP	CL-102	CL-103	CL-104	CL-104
ISIS ID:	ENCL102XXX94XD	ENCL102XXX94XX	ENCL103XXX94XX	ENCL104XXX94XX	ENCL104XXX94XX
LAB NUMBER:	2223317	2223314	2223318	2223312	2223312 D
DATE SAMPLED:	10/04/94	10/04/94	10/04/94	10/05/94	10/05/94
DATE EXTRACTED:	10/08/94	10/08/94	10/08/94	10/08/94	10/08/94
DATE ANALYZED:	10/29/94	10/29/94	10/29/94	11/04/94	11/04/94

ANALYTE	SOW-3/90 - 11	CRQL					
alpha-BHC	0.05	0.056 U	0.05 U	0.05 U	0.05 U	0.5 U	
beta-BHC	0.05	0.056 U	0.05 U	0.05 U	0.05 U	0.5 U	
delta-BHC	0.05	0.056 U	0.05 U	0.05 U	0.05 U	0.5 U	
gamma-BHC (Lindane)	0.05	0.056 U	0.05 U	0.05 U	0.05 U	0.5 U	
Heptachlor	0.05	0.056 U	0.05 U	0.05 U	0.05 U	0.5 U	
Aldrin	0.05	0.056 U	0.05 U	0.05 U	0.05 U	0.5 U	
Heptachlor Epoxide	0.05	0.056 U	0.05 U	0.05 U	0.05 U	0.5 U	
Endosulfan I	0.05	0.056 U	0.05 U	0.05 U	0.05 U	0.5 U	
Dieldrin	0.1	0.11 U	0.1 U	0.1 U	0.1 U	1.0 U	
4,4'-DDE	0.1	0.11 U	0.1 U	0.1 U	0.63 P	0.99 JP	
Endrin	0.1	0.11 U	0.1 U	0.1 U	0.1 U	1.0 U	
Endosulfan II	0.1	0.11 U	0.1 U	0.1 U	0.31	1.0 U	
4,4'-DDD	0.1	0.11 U	0.1 U	0.1 U	0.17 P	1.0 U	
Endrin Aldehyde	0.1	0.11 U	0.13 P	0.1 U	0.88 P	1.0 U	
Endosulfan Sulfate	0.1	0.11 U	0.1 U	0.1 U	0.1 U	1.0 U	
4,4'-DDT	0.1	0.11 U	0.1 U	0.1 U	0.1 U	1.0 U	
Methoxychlor	0.5	0.56 U	0.5 U	0.5 U	0.5 U	5.0 U	
Endrin Ketone	0.1	0.11 U	0.1 U	0.1 U	0.1 U	1.0 U	
alpha-Chlordane	0.05	0.056 U	0.05 U	0.05 U	0.038 JP	0.5 U	
gamma-Chlordane	0.05	0.056 U	0.05 U	0.05 U	0.085	0.5 U	
Toxaphene	5	5.6 U	5.0 U	5.0 U	5.0 U	50 U	
Aroclor-1016	1	1.1 U	1.0 U	1.0 U	1.0 U	10 U	
Aroclor-1221	2	2.2 U	2.0 U	2.0 U	2.0 U	20 U	
Aroclor-1232	1	1.1 U	1.0 U	1.0 U	1.0 U	10 U	
Aroclor-1242	1	1.1 U	1.0 U	1.0 U	1.0 U	10 U	
Aroclor-1248	1	1.1 U	1.0 U	1.0 U	1.0 U	10 U	
Aroclor-1254	1	1.1 U	1.0 U	1.0 U	1.0 U	10 U	
Aroclor-1260	1	1.1 U	1.0 U	1.0 U	1.0 U	10 U	

	1.00	1.00	1.00	1.00	10.0
Dilution Factor:	1.00	1.00	1.00	1.00	10.0
Sample Volume/Weight (ml/g):	900	1000	1000	1000	1000

Associated Method Blank:	PWB1008B	PWB1008B	PWB1008B	PBW1008A2	PBW1008A2
Associated Equipment Blank:	ENQSXX2XXX94XX	ENQSXX2XXX94XX	ENQSXX2XXX94XX	ENQSXX2XXX94XX	ENQSXX2XXX94XX
Associated Field Blank:	-	-	-	-	-

Site: SUMP LIQUIDS

U: not detected P: >25% difference between columns J: estimated

Table 1
Laboratory Report of Analysis

LOCATION:	CD-101	CD-103
ISIS ID:	ENCD101XXX94XX	ENCD103XXX94XX
LAB NUMBER:	2223303	2223304
DATE SAMPLED:	10/04/94	10/04/94
DATE EXTRACTED:	10/08/94	10/08/94
DATE ANALYZED:	11/13/94	11/13/94

ANALYTE	SOW-3/90 - 11	CRQL			
alpha-BHC	1.7	45	P	6.2	U
beta-BHC	1.7	27	U	6.2	U
delta-BHC	1.7	27	U	6.2	U
gamma-BHC (Lindane)	1.7	29	P	6.2	
Heptachlor	1.7	27	U	6.2	U
Aldrin	1.7	27	U	6.2	U
Heptachlor Epoxide	1.7	27	U	8.7	
Endosulfan I	1.7	27	U	6.2	U
Dieldrin	3.3	52	U	12	U
4,4'-DDE	3.3	52	U	12	U
Endrin	3.3	52	U	12	U
Endosulfan II	3.3	52	U	12	U
4,4'-DDD	3.3	52	U	17	P
Endrin Aldehyde	3.3	52	U	12	U
Endosulfan Sulfate	3.3	52	U	12	U
4,4'-DDT	3.3	52	U	12	U
Methoxychlor	17	270	U	62	U
Endrin Ketone	3.3	52	U	12	U
alpha-Chlordane	1.7	27	U	6.2	U
gamma-Chlordane	1.7	27	U	6.2	U
Toxaphene	170	2700	U	620	U
Aroclor-1016	33	520	U	120	U
Aroclor-1221	67	1000	U	240	U
Aroclor-1232	33	520	U	120	U
Aroclor-1242	33	520	U	120	U
Aroclor-1248	33	520	U	120	U
Aroclor-1254	33	520	U	120	U
Aroclor-1260	33	520	U	120	U

Dilution Factor:	5.00	2.00
Percent Solids:	32	55
Sample Volume\Weight (ml\g):	30.0	30.0

Associated Method Blank:	PSB1008B	PSB1008B
Associated Equipment Blank:	ENQSXX1XXX94XX	ENQSXX1XXX94XX
Associated Field Blank:	-	-

Site: SUMP SEDIMENTS

U: not detected P: >25% difference between columns

Table 1
Laboratory Report of Analysis

	LOCATION:	WT-101 DUP	WT-101	WT-102	WT-103	WT-104	WT-105	WT-106	WT-106
	ISIS ID:	ENWT101XXX94XD	ENWT101XXX94XX	ENWT102XXX94XX	ENWT103XXX94XX	ENWT104XXX94XX	ENWT105XXX94XX	ENWT106XXX94XX	ENWT106XXX94XX
	LAB NUMBER:	2223309	2223306	2223305	2223310	2223311	2223301	2223302	2223302 D
	DATE SAMPLED:	10/04/94	10/04/94	10/04/94	10/04/94	10/04/94	10/05/94	10/05/94	10/05/94
	DATE EXTRACTED:	10/25/94	10/08/94	10/08/94	10/08/94	10/08/94	10/08/94	10/08/94	10/08/94
	DATE ANALYZED:	12/02/94	11/13/94	11/13/94	11/13/94	11/13/94	11/13/94	11/14/94	11/17/94
ANALYTE	SOW-3/90 - II	CRQL							
alpha-BHC	1.7	2.0 U	170 P	18 U	29 U	14 U	1.9 U	25 P	37 U
beta-BHC	1.7	2.0 U	97 U	18 U	29 U	14 U	1.9 U	9.3 U	1000 DPE
delta-BHC	1.7	2.0 U	97 U	18 U	29 U	14 U	1.9 U	130 P	37 U
gamma-BHC (Lindane)	1.7	2.0 U	180 P	18 U	29 P	14 U	1.9 U	230 PE	37 U
Heptachlor	1.7	10 P	97 U	18 U	29 U	14 U	3.7 P	9.3 U	37 U
Aldrin	1.7	2.0 U	92 JP	28 P	48 P	14 U	1.9 U	9.3 U	37 U
Heptachlor Epoxide	1.7	2.0 U	280 P	18 U	29 U	14 U	6.2 P	9.3 U	37 U
Endosulfan I	1.7	5.2 P	250 P	18 U	96	16 P	5.9 P	79 P	37 U
Dieldrin	3.3	3.9 U	190 U	34 U	71 P	27 JP	3.4 JP	31 P	73 U
4,4'-DDE	3.3	3.9 U	670 P	50	67 P	51 P	10 P	120 P	73 U
Endrin	3.3	3.9 U	190 U	34 U	56 U	27 U	3.6 U	18 U	73 U
Endosulfan II	3.3	3.9 U	190 U	34 U	56 U	27 U	3.6 U	18 U	73 U
4,4'-DDD	3.3	3.9 U	190 U	34 U	64 P	58 P	3.6 U	18 U	73 U
Endrin Aldehyde	3.3	31 P	480 P	34 U	56 U	27 U	3.6 U	18 U	73 U
Endosulfan Sulfate	3.3	3.9 U	190 U	34 U	56 U	27 U	3.6 U	18 U	73 U
4,4'-DDT	3.3	7.1 P	190 U	34 U	62 P	27 U	5.8 P	33 P	73 U
Methoxychlor	17	20 U	970 U	180 U	290 U	140 U	19 U	93 U	370 U
Endrin Ketone	3.3	3.9 U	190 U	34 U	56 U	27 U	3.6 U	18 U	73 U
alpha-Chlordane	1.7	2.0 U	97 U	18 U	29 U	36 P	4.9 P	9.3 U	37 U
gamma-Chlordane	1.7	2.0 U	230 P	18 U	32 P	21	1.9 U	12 P	37 U
Toxaphene	170	200 U	9700 U	1800 U	2900 U	1400 U	190 U	930 U	3700 U
Aroclor-1016	33	39 U	1900 U	340 U	560 U	270 U	36 U	180 U	730 U
Aroclor-1221	67	80 U	3800 U	700 U	1100 U	560 U	74 U	370 U	1500 U
Aroclor-1232	33	39 U	1900 U	340 U	560 U	270 U	36 U	180 U	730 U
Aroclor-1242	33	39 U	1900 U	340 U	560 U	270 U	36 U	180 U	730 U
Aroclor-1248	33	39 U	1900 U	340 U	560 U	270 U	36 U	180 U	730 U
Aroclor-1254	33	39 U	1900 U	340 U	560 U	270 U	36 U	180 U	730 U
Aroclor-1260	33	39 U	1900 U	340 U	560 U	270 U	36 U	180 U	730 U
=====									
Dilution Factor:	1.00	50.0	5.00	10.0	5.00	1.00	5.00	20.0	
Percent Solids:	84	88	48	59	60	91	91	91	
Sample Volume\Weight (ml\g):	30.0	30.0	30.0	30.0	30.0	30.0	30.0	30.0	
Associated Method Blank:	PSB1025A	PSB1008B	PSB1008B	PSB1008B	PSB1008B	PSB1008B	PSB1008B	PSB1008A	
Associated Equipment Blank:	ENQSXX1XXX94XX	ENQSXX1XXX94XX	ENQSXX1XXX94XX	ENQSXX1XXX94XX	ENQSXX1XXX94XX	ENQSXX1XXX94XX	ENQSXX1XXX94XX	ENQSXX1XXX94XX	
Associated Field Blank:	-	-	-	-	-	-	-	-	

Site: WASTE-SLUDGE

U: not detected P: >25% difference between columns J: estimated

D: diluted result E: exceeds calibration range

Table 1
Laboratory Report of Analysis

LOCATION:	SS-101	SS-102 DUP	SS-102
ISIS ID:	ENSS101XXX94XX	ENSS102XXX94XD	ENSS102XXX94XX
LAB NUMBER:	2241101	2241105	2241102
DATE SAMPLED:	10/28/94	10/28/94	10/28/94
DATE EXTRACTED:	11/01/94	11/01/94	11/01/94
DATE ANALYZED:	12/08/94	12/08/94	12/08/94

ANALYTE	SOW-3/90 - II	CRQL			
alpha-BHC	1.7	9.4 U	2.3 U	2.3 U	
beta-BHC	1.7	9.4 U	2.3 U	2.3 U	
delta-BHC	1.7	9.4 U	2.3 U	2.3 U	
gamma-BHC (Lindane)	1.7	9.4 U	2.3 U	2.3 U	
Heptachlor	1.7	7.3 J	2.3 U	2.3 U	
Aldrin	1.7	9.4 U	2.3 U	2.3 U	
Heptachlor Epoxide	1.7	9.4 U	2.3 U	2.3 U	
Endosulfan I	1.7	9.4 U	2.3 U	2.3 U	
Dieldrin	3.3	18 U	4.5 U	4.5 U	
4,4'-DDE	3.3	18 U	4.5 U	4.5 U	
Endrin	3.3	18 U	4.5 U	4.5 U	
Endosulfan II	3.3	9.9 JP	5.6 P	4.5 U	
4,4'-DDD	3.3	18 U	4.5 U	4.5 U	
Endrin Aldehyde	3.3	18 U	4.5 U	4.5 U	
Endosulfan Sulfate	3.3	18 U	4.5 U	4.5 U	
4,4'-DDT	3.3	14 JP	4.5 P	4.5 U	
Methoxychlor	17	54 BJP	15 BJP	13 BJP	
Endrin Ketone	3.3	18 U	7.2 P	3.7 JP	
alpha-Chlordane	1.7	9.4 U	2.3 U	2.3 U	
gamma-Chlordane	1.7	9.4 U	2.3 U	2.3 U	
Toxaphene	170	940 U	230 U	230 U	
Aroclor-1016	33	180 U	45 U	45 U	
Aroclor-1221	67	370 U	91 U	91 U	
Aroclor-1232	33	180 U	45 U	45 U	
Aroclor-1242	33	180 U	45 U	45 U	
Aroclor-1248	33	180 U	45 U	45 U	
Aroclor-1254	33	180 U	45 U	45 U	
Aroclor-1260	33	180 U	45 U	45 U	

Dilution Factor:	5.00	1.00	1.00
Percent Solids:	90	74	74
Sample Volume\Weight (ml\g):	30.0	30.0	30.0

Associated Method Blank:	PSB1101A	PSB1101A	PSB1101A
Associated Equipment Blank:	ENQS003XXX94XX	ENQS003XXX94XX	ENQS003XXX94XX
Associated Field Blank:	-	-	-

Site: SURFACE SOILS

U: not detected J: estimated P: >25% difference between columns B: blank contamination

Table 1
Laboratory Report of Analysis

LOCATION:	BS-101	BS-102
ISIS ID:	ENBS101XX694XX	ENBS102XX694XX
LAB NUMBER:	2241106	2241107
DATE SAMPLED:	10/27/94	10/28/94
DATE EXTRACTED:	11/01/94	11/01/94
DATE ANALYZED:	12/08/94	12/08/94

ANALYTE	SOW-3/90 - 11	CRQL		
alpha-BHC	1.7	2.2	U	2.1 U
beta-BHC	1.7	2.2	U	2.1 U
delta-BHC	1.7	2.2	U	2.1 U
gamma-BHC (Lindane)	1.7	2.2	U	2.1 U
Heptachlor	1.7	2.2	U	2.1 U
Aldrin	1.7	2.2	U	2.1 U
Heptachlor Epoxide	1.7	2.2	U	2.1 U
Endosulfan I	1.7	2.2	U	2.1 U
Dieldrin	3.3	4.3	U	4.1 U
4,4'-DDE	3.3	4.3	U	4.1 U
Endrin	3.3	4.3	U	4.1 U
Endosulfan II	3.3	4.3	U	4.1 U
4,4'-DDD	3.3	4.3	U	4.1 U
Endrin Aldehyde	3.3	4.3	U	4.1 U
Endosulfan Sulfate	3.3	4.3	U	4.1 U
4,4'-DDT	3.3	4.3	U	4.1 U
Methoxychlor	17	22	U	21 U
Endrin Ketone	3.3	4.3	U	4.1 U
alpha-Chlordane	1.7	2.2	U	2.1 U
gamma-Chlordane	1.7	2.2	U	2.1 U
Toxaphene	170	220	U	210 U
Aroclor-1016	33	43	U	41 U
Aroclor-1221	67	87	U	83 U
Aroclor-1232	33	43	U	41 U
Aroclor-1242	33	43	U	41 U
Aroclor-1248	33	43	U	41 U
Aroclor-1254	33	43	U	41 U
Aroclor-1260	33	43	U	35 JP

Dilution Factor:	1.00	1.00
Percent Solids:	77	81
Sample Volume\Weight (mL\g):	30.0	30.0

Associated Method Blank:	PSB1101A	PSB1101A
Associated Equipment Blank:	ENQS003XXX94XX	ENQS003XXX94XX
Associated Field Blank:	-	-

Site: SOIL BORINGS

U: not detected J: estimated P: >25% difference between columns

Table 1
Laboratory Report of Analysis

LOCATION:	MW-101 DUP	MW-101	MW-102	MW-103
ISIS ID:	ENMW101XXX94XD	ENMW101XXX94XX	ENMW102XXX94XX	ENMW103XXX94XX
LAB NUMBER:	2263804	2263801	2263805	2263806
DATE SAMPLED:	11/30/94	11/30/94	11/30/94	11/30/94
DATE EXTRACTED:	12/03/94	12/03/94	12/03/94	12/03/94
DATE ANALYZED:	12/12/94	12/12/94	12/12/94	12/12/94

ANALYTE	SOW-3/90 - II	CRQL				
alpha-BHC	0.05	0.05 U	0.05 U	0.05 U	0.05 U	0.05 U
beta-BHC	0.05	0.05 U	0.05 U	0.05 U	0.05 U	0.05 U
delta-BHC	0.05	0.05 U	0.05 U	0.05 U	0.05 U	0.05 U
gamma-BHC (Lindane)	0.05	0.05 U	0.05 U	0.05 U	0.05 U	0.05 U
Heptachlor	0.05	0.05 U	0.05 U	0.05 U	0.05 U	0.05 U
Aldrin	0.05	0.05 U	0.05 U	0.05 U	0.05 U	0.05 U
Heptachlor Epoxide	0.05	0.05 U	0.05 U	0.05 U	0.05 U	0.05 U
Endosulfan I	0.05	0.05 U	0.05 U	0.05 U	0.05 U	0.05 U
Dieldrin	0.1	0.1 U	0.1 U	0.1 U	0.1 U	0.1 U
4,4'-DDE	0.1	0.1 U	0.1 U	0.1 U	0.1 U	0.1 U
Endrin	0.1	0.1 U	0.1 U	0.1 U	0.1 U	0.1 U
Endosulfan II	0.1	0.1 U	0.1 U	0.1 U	0.1 U	0.1 U
4,4'-DDD	0.1	0.1 U	0.1 U	0.1 U	0.1 U	0.1 U
Endrin Aldehyde	0.1	0.1 U	0.1 U	0.1 U	0.1 U	0.1 U
Endosulfan Sulfate	0.1	0.1 U	0.1 U	0.1 U	0.1 U	0.1 U
4,4'-DDT	0.1	0.1 U	0.1 U	0.1 U	0.1 U	0.1 U
Methoxychlor	0.5	0.5 U	0.5 U	0.5 U	0.5 U	0.5 U
Endrin Ketone	0.1	0.1 U	0.1 U	0.1 U	0.1 U	0.1 U
alpha-Chlordane	0.05	0.05 U	0.05 U	0.05 U	0.05 U	0.05 U
gamma-Chlordane	0.05	0.05 U	0.05 U	0.05 U	0.05 U	0.05 U
Toxaphene	5	5.0 U	5.0 U	5.0 U	5.0 U	5.0 U
Aroclor-1016	1	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
Aroclor-1221	2	2.0 U	2.0 U	2.0 U	2.0 U	2.0 U
Aroclor-1232	1	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
Aroclor-1242	1	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
Aroclor-1248	1	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
Aroclor-1254	1	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
Aroclor-1260	1	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U

Dilution Factor:	1.00	1.00	1.00	1.00
Sample Volume/Weight (ml\g):	1000	1000	1000	1000

Associated Method Blank:	PWB1203A	PWB1203B	PWB1203A	PWB1203A
Associated Equipment Blank:	ENQS004XXX94XX	ENQS004XXX94XX	ENQS004XXX94XX	ENQS004XXX94XX
Associated Field Blank:	-	-	-	-

Site: MONITORING WELLS
U: not detected

Table 1
Laboratory Report of Analysis

LOCATION:	QS-003	QS-004	QS-XX1	QS-2
ISIS ID:	ENQS003XXX94XX	ENQS004XXX94XX	ENQSXX1XXX94XX	ENQSXX2XXX94XX
LAB NUMBER:	241108	263807	223319	223313
DATE SAMPLED:	10/28/94	11/30/94	10/04/94	10/04/94

ANALYTE	SOW-3/90 - 11	CRDL				
Aluminum	200		57.0 U	57.0 U	57.0 U	57.0 U
Antimony	60		38.0 U	38.0 U	38.0 U	38.0 U
Arsenic	10		5.0 UN	5.0 U	5.0 U	5.0 U
Barium	200		11.0 U	11.0 U	11.0 U	11.0 U
Beryllium	5		2.0 U	2.0 U	2.0 U	2.0 U
Cadmium	5		2.0 U	2.0 U	2.0 U	2.0 U
Calcium	5000		1390 U*	1390 U	1390 U	1390 U
Chromium	10		5.0 U	5.0 U	5.0 U*	5.0 U*
Cobalt	50		6.0 U	6.0 U	6.0 U	6.0 U
Copper	25		5.0 U	5.0 U	5.0 U	5.0 U
Iron	100		28.4 B	16.0 U*	28.9 B	16.0 U
Lead	3		3.0 U*	3.0 UN	3.0 U*	6.1 *
Magnesium	5000		1550 U	1550 U	1550 U	1550 U
Manganese	15		2.0 U*	2.0 U	2.0 UN	2.0 UN
Mercury	0.2		0.20 U	0.20 U	0.20 U*	0.20 U*
Nickel	40		26.0 U	26.0 U	26.0 U	26.0 U
Potassium	5000		840 U	840 U	840 U	840 U
Selenium	5		5.0 UN	5.0 U	5.0 UN	5.0 UN
Silver	10		5.0 U	5.0 U	5.0 U	5.0 U
Sodium	5000		463 U	479 B	463 U	463 U
Thallium	10		5.0 U	5.0 U	5.0 U	5.0 U
Vanadium	50		17.0 U	17.0 U	17.0 U	17.0 U
Zinc	20		9.1 B	5.0 U	5.4 B	5.0 U
Cyanide	10		10.0 UN	10.0 U	10.0 U	10.0 U

Associated Method Blank:	PBENRX2	PBENRX3	PBENRX1	PBENRX1
Associated Equipment Blank:	-	-	-	-
Associated Field Blank:	-	-	-	-

Site: EQUIPMENT RINSATE

U: not detected N: spike recovery not met *: duplicate analysis not met

B: less than CRDL W: post digestion spike not met

Table 1
Laboratory Report of Analysis

LOCATION: QS-003
ISIS ID: ENQS003XXX94XX
LAB NUMBER: D241108
DATE SAMPLED: 10/28/94

ANALYTE	SOW-3/90 - II	CRDL	
Aluminum	200	57.0	U
Antimony	60	38.0	U
Arsenic	10	5.0	UN
Barium	200	11.0	U
Beryllium	5	2.0	U
Cadmium	5	2.0	U
Calcium	5000	1390	U*
Chromium	10	5.0	U
Cobalt	50	6.0	U
Copper	25	5.0	U
Iron	100	16.0	U
Lead	3	3.0	U*
Magnesium	5000	1550	U
Manganese	15	2.0	U*
Mercury	0.2	0.20	U
Nickel	40	26.0	U
Potassium	5000	840	U
Selenium	5	5.0	UN
Silver	10	5.0	U
Sodium	5000	463	U
Thallium	10	NR	
Vanadium	50	17.0	U
Zinc	20	5.0	U
Cyanide	10	NR	

Associated Method Blank: PBENRX2
Associated Equipment Blank: -
Associated Field Blank: -

Site: EQUIPMENT RINSATE

U: not detected N: spike recovery not met *: duplicate analysis not met

Table 1
Laboratory Report of Analysis

LOCATION:	CL-102 DUP	CL-102	CL-103	CL-104
ISIS ID:	ENCL102XXX94XD	ENCL102XXX94XX	ENCL103XXX94XX	ENCL104XXX94XX
LAB NUMBER:	223317	223314	223318	223312
DATE SAMPLED:	10/04/94	10/04/94	10/04/94	10/05/94

ANALYTE	SOW-3/90 - 11	CRDL				
Aluminum	200	9040	9520	366	27000	
Antimony	60	38.0 U	38.0 U	38.0 U	38.0 U	
Arsenic	10	5.3 B	5.0 UW	5.0 U	54.2	
Barium	200	180 B	203	39.6 B	1320	
Beryllium	5	2.0 U	2.0 U	2.0 U	2.0 U	
Cadmium	5	29.2	32.7	2.1 B	122	
Calcium	5000	53800	58400	68800	580000	
Chromium	10	137 *	147 *	9.7 B*	1870 *	
Cobalt	50	30.7 B	38.0 B	21.4 B	375	
Copper	25	1820	1970	120	4930	
Iron	100	70000	91200	20000	265000	
Lead	3	1960	1770 *	94.3 S*	5510	
Magnesium	5000	10600	12100	18300	57000	
Manganese	15	1300 W	1610 W	2710 W	2330 W	
Mercury	0.2	1.6 *	2.0 *	0.20 U*	545 *	
Nickel	40	155	196	48.7	3500	
Potassium	5000	4600 B	5350	13200	30800	
Selenium	5	5.0 UWN	25.0 UWN	5.0 UWN	10.0 UN	
Silver	10	5.0 U	5.0 U	5.0 U	23.1	
Sodium	5000	5840	5870	38300	2030000	
Thallium	10	5.0 U	5.0 U	5.0 UW	5.0 UW	
Vanadium	50	67.3	76.5	17.0 U	124	
Zinc	20	7190	8540	337	24500	
Cyanide	10	10.0 U	10.0 U	10.0 U	10.0 U	

Associated Method Blank:	PBENRX1	PBENRX1	PBENRX1	PBENRX1
Associated Equipment Blank:	ENQSXX2XXX94XX	ENQSXX2XXX94XX	ENQSXX2XXX94XX	ENQSXX2XXX94XX
Associated Field Blank:	-	-	-	-

Site: SUMP LIQUIDS

U: not detected B: less than CRDL *: duplicate analysis not met

N: spike recovery not met W: post digestion spike not met

Table 1
Laboratory Report of Analysis

LOCATION:	CD-101	CD-103
ISIS ID:	ENCD101XXX94XX	ENCD103XXX94XX
LAB NUMBER:	223303	223304
DATE SAMPLED:	10/04/94	10/04/94

ANALYTE	SOW-3/90 - 11	CRDL		
Aluminum	40	1750	4270	
Antimony	12	23.1 UN	16.1 BN	
Arsenic	2	6.6 N	6.8 SN	
Barium	40	142	186	
Beryllium	1	1.2 U	0.67 U	
Cadmium	1	11.6	10.2	
Calcium	1000	4460	17800	
Chromium	2	78.4 *	254 *	
Cobalt	10	22.7 B*	45.1 *	
Copper	5	181	650	
Iron	20	66400 *	237000 *	
Lead	0.6	246 S*	213 *	
Magnesium	1000	943 U	3550	
Manganese	3	974 *	1290 *	
Mercury	0.1	1.3	1.0	
Nickel	8	167 *	406 *	
Potassium	1000	512 U	329 B	
Selenium	1	3.0 UWN*	1.7 UN*	
Silver	2	3.0 UN*	1.7 UN*	
Sodium	1000	2050 B	360 B	
Thallium	2	3.0 UN	1.7 UN	
Vanadium	10	16.4 BN	59.3 N	
Zinc	4	1140 *	1140 *	
Cyanide	1	1.5 U	0.94 U	

Percent Solids: 32 55

Associated Method Blank:	PBENRX1	PBENRX1
Associated Equipment Blank:	ENQSXX1XXX94XX	ENQSXX1XXX94XX
Associated Field Blank:	-	-

Site: SUMP SEDIMENTS

U: not detected N: spike recovery not met *: duplicate analysis not met

B: less than CRDL S: method of standard additions W: post digestion spike not met

Table 1
Laboratory Report of Analysis

LOCATION:	WT-101 DUP	WT-101	WT-102	WT-103	WT-104	WT-105	WT-106
ISIS ID:	ENWT101XXX94XD	ENWT101XXX94XX	ENWT102XXX94XX	ENWT103XXX94XX	ENWT104XXX94XX	ENWT105XXX94XX	ENWT106XXX94XX
LAB NUMBER:	223309	223306	223305	223310	223311	223301	223302
DATE SAMPLED:	10/04/94	10/04/94	10/04/94	10/04/94	10/04/94	10/05/94	10/05/94

ANALYTE	SOW-3/90 - 11	CRDL							
Aluminum	40	6370		3910		6380		1870	
Antimony	12	8.7 UN		8.3 UN		13.6 BN		12.8 UN	
Arsenic	2	11.3 N		9.6 N		35.6 N		21.7 SN	
Barium	40	419		345		184		110	
Beryllium	1	0.83 B		0.69 B		0.69 U		0.58 U	
Cadmium	1	36.8		21.4		9.1		18.6	
Calcium	1000	7110		6720		68000		21300	
Chromium	2	3730 *		2790 *		346 *		102 *	
Cobalt	10	861 *		912 *		29.7 *		34.8 *	
Copper	5	2770		1800		379		553	
Iron	20	56900 *		133000 *		112000 *		608000 *	
Lead	0.6	1810 *		1850 *		363 *		275 *	
Magnesium	1000	1650		1330		5530		2900	
Manganese	3	501 *		916 *		891 *		2830 *	
Mercury	0.1	1.1		1.1		52.9		1.8	
Nickel	8	11400 *		9530 *		209 *		129 *	
Potassium	1000	1490		1240		1160 B		2460	
Selenium	1	5.5 UN*		2.1 SN*		1.9 UWN*		1.5 UN*	
Silver	2	16.6 N*		20.5 N*		1.7 UN*		1.7 UN*	
Sodium	1000	6060		3680		984 B		3050	
Thallium	2	1.1 UWN		1.0 UWN		1.9 UN		1.5 UN	
Vanadium	10	30.2 N		84.3 N		45.3 N		58.4 N	
Zinc	4	2300 *		2090 *		1910 *		1370 *	
Cyanide	1	0.52 U		2.7		0.99 U		20.7	
Percent Solids:			84	88	48	59	60	91	91

Associated Method Blank:	PBENRX1	PBENRX1	PBENRX1	PBENRX1	PBENRX1	PBENRX1	PBENRX1
Associated Equipment Blank:	ENQSXX1XXX94XX	ENQSXX1XXX94XX	ENQSXX1XXX94XX	ENQSXX1XXX94XX	ENQSXX1XXX94XX	ENQSXX1XXX94XX	ENQSXX1XXX94XX
Associated Field Blank:	-	-	-	-	-	-	-

Site: WASTE-SLUDGE

U: not detected N: spike recovery not met B: less than CRDL

*: duplicate analysis not met W: post digestin spike not met S: method of standard additions

Table 1
Laboratory Report of Analysis

LOCATION:	SS-101	SS-102 DUP	SS-102
ISIS ID:	ENSS101XXX94XX	ENSS102XXX94XD	ENSS102XXX94XX
LAB NUMBER:	241101	241105	241102
DATE SAMPLED:	10/28/94	10/28/94	10/28/94

ANALYTE	SOW-3/90 - II	CRDL			
Aluminum	40	1650		9860	10200
Antimony	12	8.0 U		9.2 U	10.1 U
Arsenic	2	4.4 N		35.1 SN	40.5 SN
Barium	40	27.3 B		142	178
Beryllium	1	0.42 U		4.7	4.0
Cadmium	1	1.8		1.1 B	2.3
Calcium	1000	212000 *		34900 *	35900 *
Chromium	2	11.4		10.1	14.3
Cobalt	10	6.1 B		10.9 B	12.1 B
Copper	5	67.9		77.9	69.4
Iron	20	7870		12800	18200
Lead	0.6	95.1 N		136 N	160 S*
Magnesium	1000	18300		7060	6400
Manganese	3	495 *		380 *	284 *
Mercury	0.1	0.40		0.61	0.88
Nickel	8	24.5		25.9	30.8
Potassium	1000	470 B		521 B	1010 B
Selenium	1	1.1 UWN		1.3 UWN	1.3 UWN
Silver	2	1.0 U		1.2 U	1.3 U
Sodium	1000	106 B		237 B	190 B
Thallium	2	1.1 UW		1.3 U	1.3 U
Vanadium	10	11.7		34.4	35.4
Zinc	4	653		2100	1940
Cyanide	1	0.53 UN		0.57 UN	0.58 UN
=====					
Percent Solids:		90	74	74	

Associated Method Blank: PBENRX2 PBENRX2 PBENRX2
Associated Equipment Blank: ENQS003XXX94XX ENQS003XXX94XX ENQS003XXX94XX
Associated Field Blank: - - -

Site: SURFACE SOILS

U: not detected N: spike recovery not met B: less than CRDL

*: duplicate analysis not met W: post digestion spike not met S: method of standard additions

Table 1
Laboratory Report of Analysis

LOCATION:	BS-101	BS-102
ISIS ID:	ENBS101XX694XX	ENBS102XX694XX
LAB NUMBER:	241106	241107
DATE SAMPLED:	10/27/94	10/28/94

ANALYTE	SOW-3/90 - II	CRDL			
Aluminum	40	7130		16100	
Antimony	12	9.2 U		9.4 B	
Arsenic	2	5.1 N		4.6 N	
Barium	40	57.0		186	
Beryllium	1	0.48 U		0.93 B	
Cadmium	1	0.48 U		0.47 U	
Calcium	1000	12200 *		19700 *	
Chromium	2	8.7		20.8	
Cobalt	10	6.4 B		10.2 B	
Copper	5	49.5		25.9	
Iron	20	15600		31100	
Lead	0.6	50.5 N		26.0 S*	
Magnesium	1000	4780		8670	
Manganese	3	270 *		485 *	
Mercury	0.1	0.13 U		0.12 U	
Nickel	8	12.9		26.3	
Potassium	1000	982 B		2470	
Selenium	1	1.1 UWN		1.2 UN	
Silver	2	1.2 U		1.2 U	
Sodium	1000	1330		295 B	
Thallium	2	1.1 UW		1.2 U	
Vanadium	10	19.9		42.6	
Zinc	4	65.4		117	
Cyanide	1	0.67 UN		0.52 UN	
=====					
Percent Solids:		77		81	

Associated Method Blank:	PBENRX2	PBENRX2
Associated Equipment Blank:	ENQS003XXX94XX	ENQS003XXX94XX
Associated Field Blank:	-	-

Site: SOIL BORINGS

U: not detected N: spike recovery not met *: duplicate analysis not met

B: less than CRDL S: method of standard additions W: post digestion spike not met

Table 1
Laboratory Report of Analysis

LOCATION:	MW-101 DUP	MW-101	MW-102	MW-103
ISIS ID:	ENMW101XXX94XD	ENMW101XXX94XX	ENMW102XXX94XX	ENMW103XXX94XX
LAB NUMBER:	263804	263801	263805	263806
DATE SAMPLED:	11/30/94	11/30/94	11/30/94	11/30/94

ANALYTE	SOW-3/90 - 11	CRDL				

Aluminum	200	2390		250	315	57.0 U
Antimony	60	52.9 B		38.0 U	38.0 U	47.1 B
Arsenic	10	5.0 U		5.0 U	5.0 U	5.0 U
Barium	200	88.7 B		82.5 B	80.4 B	400
Beryllium	5	2.0 U		2.0 U	2.0 U	2.0 U
Cadmium	5	2.0 U		2.0 U	2.0 U	2.0 U
Calcium	5000	136000		128000	212000	137000
Chromium	10	7.2 B		5.0 U	5.0 U	5.0 U
Cobalt	50	6.0 U		6.0 U	6.0 U	6.0 U
Copper	25	11.8 B		5.0 U	5.4 B	7.5 B
Iron	100	7760 *		6240 *	1850 *	23600 *
Lead	3	11.2		3.0 U	15.7	8.5
Magnesium	5000	54600		53100	41000	36200
Manganese	15	412		339	2990	5770
Mercury	0.2	0.20 U		0.20 U	0.20 U	0.20 U
Nickel	40	26.0 U		26.0 U	26.0 U	26.0 U
Potassium	5000	14200		14200	49200	14200
Selenium	5	5.0 U		5.0 U	5.0 UW	5.0 U
Silver	10	5.0 U		5.0 U	5.0 U	5.0 U
Sodium	5000	107000		110000	208000	63300
Thallium	10	5.0 U		5.0 U	5.0 U	5.0 U
Vanadium	50	17.0 U		17.0 U	17.0 U	17.0 U
Zinc	20	13.6 B		5.0 U	24.7	17.5 B
Cyanide	10	10.0 U		10.0 U	10.0 U	10.0 U
=====						

Associated Method Blank:	PBENRX3	PBENRX3	PBENRX3	PBENRX3
Associated Equipment Blank:	ENQS004XXX94XX	ENQS004XXX94XX	ENQS004XXX94XX	ENQS004XXX94XX
Associated Field Blank:	-	-	-	-

Site: MONITORING WELLS

U: not detected B: less than CRDL *: duplicate analysis not met W: post digestion spike not met

Table 1
Laboratory Report of Analysis

LOCATION:	QS-1	QS-2
ISIS ID:	ENQSXX1XXX94XX	ENQSXX2XXX94XX
LAB NUMBER:	D223319	D223313
DATE SAMPLED:	10/04/94	10/04/94

ANALYTE	RL		
arsenic	5.0	5.0 U	5.0 U
barium	11.0	11.0 U	11.0 U
cadmium	2.0	2.0 U	2.0 U
chromium	5.0	5.0 U*	5.0 U*
lead	3.0	3.0 UW*	3.0 U*
mercury	0.20	0.20 U*	0.20 U*
selenium	5.0	5.0 UN	5.0 UN
silver	5.0	5.0 U	5.0 U

=====

Associated Method Blank:	EPENRX1	EPENRX1
Associated Equipment Blank:	-	-
Associated Field Blank:	-	-

SITE: EQUIPMENT RINSATES

Note: Inorganic Data - EPTOX Metals

U: not detected *: duplicate analysis not met W: post digestion spike not met N: spike recovery not met

Table 1
Laboratory Report of Analysis

LOCATION:	CL-103	CL-104
ISIS ID:	ENCL103XXX94XX	ENCL104XXX94XX
LAB NUMBER:	D223318	D223312
DATE SAMPLED:	10/04/94	10/05/94

ANALYTE	RL		
arsenic	52.0/5.0	5.0 U	10.0 UW
barium	11.0	41.1 B	98.6 B
cadmium	2.0	3.8 B	2.0 U
chromium	5.0	7.9 B*	5.0 U*
lead	26.0/3.0	94.3 S*	146 *
mercury	0.20	0.20 U*	0.57 *
selenium	90.0/5.0	5.0 UWN	10.0 UWN
silver	5.0	5.0 U	5.0 U

=====
Associated Method Blank: EPENRX1 EPENRX1
Associated Equipment Blank: ENQSXX2XXX94XX ENQSXX2XXX94XX
Associated Field Blank: - -

SITE: SUMP LIQUIDS

Note: Inorganic Data - EPTOX Metals

U: not detected B: less than CRDL *: duplicate analysis not met S: method of standard additions
W: post digestion spike not met N: spike recovery not met

Table 1
Laboratory Report of Analysis

LOCATION:	CD-101	CD-103
ISIS ID:	ENCD101XXX94XX	ENCD103XXX94XX
LAB NUMBER:	E223303	E223304
DATE SAMPLED:	10/04/94	10/04/94

ANALYTE	RL		
arsenic	52.0/5.0	52.0 UN	52.0 UN
barium	11.0	327 N	208 N
cadmium	2.0	2.0 U	2.0 U
chromium	5.0	5.0 U*	5.0 U*
lead	26.0/3.0	26.0 U*	26.0 U*
mercury	0.20	0.20 U	0.20 U
selenium	90.0/5.0	90.0 U	90.0 U
silver	5.0	5.0 U*	5.0 U*

=====

Associated Method Blank:	EPENRX1	EPENRX1
Associated Equipment Blank:	ENQSXX1XXX94XX	ENQSXX1XXX94XX
Associated Field Blank:	-	-

SITE: SUMP SEDIMENTS

Note: Inorganic Data - EPTOX Metals

U: not detected N: spike recovery not met *: duplicate analysis not met

Table 1
Laboratory Report of Analysis

ANALYTE	LOCATION: WT-101 DUP WT-101 WT-102 WT-103 WT-104 WT-105 WT-106 ISIS ID: ENWT101XXX94XD ENWT101XXX94XX ENWT102XXX94XX ENWT103XXX94XX ENWT104XXX94XX ENWT105XXX94XX ENWT106XXX94XX LAB NUMBER: E223309 E223306 E223305 E223310 E223311 E223301 E223302 DATE SAMPLED: 10/04/94 10/04/94 10/04/94 10/04/94 10/04/94 10/05/94 10/05/94							
	RL							
arsenic	52.0/5.0	52.0 UN	52.0 UN	52.0 UN	52.0 UN	52.0 UN	52.0 UN	52.0 UN
barium	11.0	469 N	317 N	475 N	427 N	229 N	411 N	401 N
cadmium	2.0	1100	1100	3.7 B	22.5	2.0 U	17.0 U	10.9
chromium	5.0	84.3 *	51.0 *	5.9 B*	9.5 B*	5.0 U*	53.6 *	25.0 *
lead	26.0/3.0	4820 *	11400 *	26.0 U*	58.1 *	26.0 U*	45.2 *	60.8 *
mercury	0.20	0.25	0.34	0.20 U	0.23	0.20 U	0.52	1.2
selenium	90.0/5.0	90.0 U	90.0 U	90.0 U	90.0 U	90.0 U	90.0 U	90.0 U
silver	5.0	37.6 *	33.4 *	5.0 U*	5.0 U*	5.0 U*	5.0 U*	5.0 U*

=====

Associated Method Blank:	EPENRX1	EPENRX1	EPENRX1	EPENRX1	EPENRX1	EPENRX1	EPENRX1
Associated Equipment Blank:	ENQSXX1XXX94XX	ENQSXX1XXX94XX	ENQSXX1XXX94XX	ENQSXX1XXX94XX	ENQSXX1XXX94XX	ENQSXX1XXX94XX	ENQSXX1XXX94XX
Associated Field Blank:	-	-	-	-	-	-	-

SITE: WASTE-SLUDGE

Note: Inorganic Data - EPTOX Metals

U: not detected N: spike recovery not met B: less than CRDL *: duplicate analysis not met

Table 1
Laboratory Report of Analysis

LOCATION:	SS-101	SS-102 DUP	SS-102
ISIS ID:	ENSS101XXX94XX	ENSS102XXX94XD	ENSS102XXX94XX
LAB NUMBER:	E241101	E241105	E241102
DATE SAMPLED:	10/28/94	10/28/94	10/28/94

ANALYTE	RL			
arsenic	52.0	52.0 U	52.0 U	52.0 U
barium	11.0	507	1000	1080
cadmium	2.0	2.0 U	15.8	15.8
chromium	5.0	5.0 U	6.0 B	5.0 U
lead	26.0	26.0 U	33.7	33.7
mercury	0.2	0.2 U	0.2 U	0.2 U
selenium	90.0	90.0 U	90.0 U	90.0 U
silver	5.0	5.0 U	5.0 U	5.0 U

=====

Associated Method Blank:	EPENRX2	EPENRX2	EPENRX2
Associated Equipment Blank:	-	-	-
Associated Field Blank:	-	-	-

SITE: SURFACE SOILS

Note: Inorganic Data - EPTOX Metals

U: not detected B: less than CRDL

PROJECT: NYSDEC-PSA-14 ENRX Inc. Site

Miscellaneous Aqueous Analysis

14-Feb-95

Table 1
Laboratory Report of Analysis

	LOCATION: QS-003	QS-1	QS-2
ISIS ID:	ENQS003XXX94XX	ENQSXX1XXX94XX	ENQSXX2XXX94XX
LAB NUMBER:	2241108	2223319	2223313
DATE SAMPLED:	10/28/94	10/04/94	10/04/94
DATE ANALYZED:	11/14/94	10/12/94	10/12/94

ANALYTE	RL			
Corrosivity, inch/Year	0.01	0.01 U	0.01 U	0.01 U
Ignitability, Degrees F		>212	>212	>212
Cyanide, Reactive, ppm	1	1 U	1 U	1 U
Sulfide, Reactive, ppm	1	1 U	1 U	1 U

=====

Associated Method Blank:	WCENRX2	WCENRX1	WCENRX1
Associated Equipment Blank:	-	-	-
Associated Field Blank:	-	-	-

SITE: EQUIPMENT RINSATES

U: not detected

Table 1
Laboratory Report of Analysis

LOCATION:	CL-103	CL-104
ISIS ID:	ENCL103XXX94XX	ENCL104XXX94XX
LAB NUMBER:	2223318	2223312
DATE SAMPLED:	10/04/94	10/05/94
DATE ANALYZED:	10/12/94	10/12/94

ANALYTE	RL		
Corrosivity, inch/Year	0.01	0.01 U	0.01 U
Ignitability, Degrees F		>212	>212
Cyanide, Reactive, ppm	1	1 U	1 U
Sulfide, Reactive, ppm	1	1 U	1 U

=====

Associated Method Blank:	WCENRX1	WCENRX1
Associated Equipment Blank:	ENQSXX2XXX94XX	ENQSXX2XXX94XX
Associated Field Blank:	-	-

SITE: SUMP LIQUIDS
U: not detected

Table 1
Laboratory Report of Analysis

	WT-101 DUP	WT-101	WT-102	WT-103	WT-104	WT-105	WT-106
LOCATION:	WT-101 DUP	WT-101	WT-102	WT-103	WT-104	WT-105	WT-106
ISIS ID:	ENWT101XXX94XD	ENWT101XXX94XX	ENWT102XXX94XX	ENWT103XXX94XX	ENWT104XXX94XX	ENWT105XXX94XX	ENWT106XXX94XX
LAB NUMBER:	2223309	2223306	2223305	2223310	2223311	2223301	2223302
DATE SAMPLED:	10/04/94	10/04/94	10/04/94	10/04/94	10/04/94	10/05/94	10/05/94
DATE ANALYZED:	10/12/94	10/12/94	10/12/94	10/12/94	10/12/94	10/12/94	10/12/94
ANALYTE	RL						
Corrosivity, inch/Year	0.01	0.01 U	0.01 U	0.01 U	0.01 U	0.01 U	0.01 U
Ignitability, Degrees F		>212	>212	>212	>212	>212	>212
Cyanide, Reactive, ppm	1	1 U	1 U	1 U	1 U	1 U	1 U
Sulfide, Reactive, ppm	1	1 U	1 U	1 U	1 U	1 U	1 U

```

=====
Associated Method Blank:  WCENRX1      WCENRX1      WCENRX1      WCENRX1      WCENRX1      WCENRX1      WCENRX1
Associated Equipment Blank: ENQSXX1XXX94XX ENQSXX1XXX94XX ENQSXX1XXX94XX ENQSXX1XXX94XX ENQSXX1XXX94XX ENQSXX1XXX94XX ENQSXX1XXX94XX
Associated Field Blank:    -            -            -            -            -            -            -

```

SITE: WASTE-SLUDGE

U: not detected

Table 1
Laboratory Report of Analysis

LOCATION:	CD-101	CD-103
ISIS ID:	ENCD101XXX94XX	ENCD103XXX94XX
LAB NUMBER:	2223303	2223304
DATE SAMPLED:	10/04/94	10/04/94
DATE ANALYZED:	10/12/94	10/12/94

ANALYTE	RL		
Corrosivity, inch/Year	0.01	0.01 U	0.01 U
Ignitability, Degrees F		>212	>212
Cyanide, Reactive, ppm	1	1 U	1 U
Sulfide, Reactive, ppm	1	1 U	1 U

=====

Associated Method Blank:	WCENRX1	WCENRX1
Associated Equipment Blank:	ENQSXX1XXX94XX	ENQSXX1XXX94XX
Associated Field Blank:	-	-

SITE: SUMP SEDIMENTS
U: not detected

Table 1
Laboratory Report of Analysis

ANALYTE	RL			
LOCATION:	SS-101	SS-102 DUP	SS-102	
ISIS ID:	ENSS101XXX94XX	ENSS102XXX94XD	ENSS102XXX94XX	
LAB NUMBER:	2241101	2241105	2241102	
DATE SAMPLED:	10/28/94	10/28/94	10/28/94	
DATE ANALYZED:	11/14/94	11/14/94	11/14/94	
Corrosivity, inch/Year	0.01	0.01 U	0.01 U	0.01 U
Ignitability, Degrees F		>212	>212	>212
Cyanide, Reactive, ppm	1	1 U	1 U	1 U
Sulfide, Reactive, ppm	1	1 U	1 U	1 U

=====

Associated Method Blank:	WCENRX2	WCENRX2	WCENRX2
Associated Equipment Blank:	ENQS003XXX94XX	ENQS003XXX94XX	ENQS003XXX94XX
Associated Field Blank:	-	-	-

SITE: SURFACE SOILS

U: not detected

Table 2
Validation / Summary Table

LOCATION:	CL-102 DUP	CL-102	CL-103	CL-104
ISIS ID:	ENCL102XXX94XD	ENCL102XXX94XX	ENCL103XXX94XX	ENCL104XXX94XX
LAB NUMBER:	2223317	2223314	2223318	2223312
DATE SAMPLED:	10/04/94	10/04/94	10/04/94	10/05/94
DATE ANALYZED:	10/13/94	10/13/94	10/13/94	10/13/94

ANALYTE	SOW-3/90 - II	CRQL				
Chloromethane	10	10 U	10 U	10 U	200 U	
Bromomethane	10	10 U	10 U	10 U	200 U	
Vinyl Chloride	10	34 J	80 J	10 U	230	
Chloroethane	10	10 U	10 U	10 U	200 U	
Methylene Chloride	10	18 U	17 U	16 U	830 U	
Acetone	10	16 UJ	10 U	10 U	200 U	
Carbon Disulfide	10	10 U	10 U	10 U	200 U	
1,1-Dichloroethene	10	10 U	10 U	10 U	200 U	
1,1-Dichloroethane	10	16	15	10 U	12000	
1,2-Dichloroethene (total)	10	480	480	10 U	13000	
Chloroform	10	10 U	10 U	10 U	200 U	
1,2-Dichloroethane	10	10 U	10 U	10 U	200 U	
2-Butanone	10	10 U	10 U	10 U	200 U	
1,1,1-Trichloroethane	10	190	170	10 U	57000	
Carbon Tetrachloride	10	10 U	10 U	10 U	200 U	
Bromodichloromethane	10	10 U	10 U	10 U	200 U	
1,2-Dichloropropane	10	10 U	10 U	10 U	200 U	
cis-1,3-Dichloropropene	10	10 U	10 U	10 U	200 U	
Trichloroethene	10	10 J	6 J	10 U	33000	
Dibromochloromethane	10	10 U	10 U	10 U	200 U	
1,1,2-Trichloroethane	10	10 U	10 U	10 U	200 U	
Benzene	10	10 U	10 U	10 U	200 U	
trans-1,3-Dichloropropene	10	10 U	10 U	10 U	200 U	
Bromoform	10	10 U	10 U	10 U	200 U	
4-Methyl-2-Pentanone	10	10 U	10 U	10 U	200 U	
2-Hexanone	10	10 U	10 U	10 U	200 U	
Tetrachloroethene	10	4 J	2 J	10 U	8000 J	
1,1,2,2-Tetrachloroethane	10	10 U	10 U	10 U	200 U	
Toluene	10	28 J	13 J	10 U	510	
Chlorobenzene	10	10 U	10 U	10 U	200 U	
Ethylbenzene	10	8 J	3 J	10 U	43 J	
Styrene	10	10 U	10 U	10 U	200 U	
Total Xylenes	10	30 J	10 J	10 U	210	

Dilution Factor:	1.00	1.00	1.00	20.0
Sample Volume\Weight (ml\g):	5.00	5.00	5.00	5.00
Associated Method Blank:	N9594.D	N9594.D	N9594.D	N9594.D
Associated Equipment Blank:	ENQSXX2XXX94XX	ENQSXX2XXX94XX	ENQSXX2XXX94XX	ENQSXX2XXX94XX
Associated Field Blank:	-	-	-	-
Associated Trip Blank:	ENQT001XXX94XX	ENQT001XXX94XX	ENQT001XXX94XX	ENQT001XXX94XX

Site: SUMP LIQUIDS

U: not detected J: estimated

Table 2
Validation / Summary Table

		LOCATION:	CD-101	CD-103
		ISIS ID:	ENCD101XXX94XX	ENCD103XXX94XX
		LAB NUMBER:	2223303	2223304
		DATE SAMPLED:	10/04/94	10/04/94
		DATE ANALYZED:	10/13/94	10/13/94
ANALYTE	SOW-3/90 - II	CRQL		
Chloromethane	10	160 U	18 U	
Bromomethane	10	160 U	18 U	
Vinyl Chloride	10	390	18 U	
Chloroethane	10	160 U	18 U	
Methylene Chloride	10	280 U	32 U	
Acetone	10	960 J	150 J	
Carbon Disulfide	10	160 U	18 U	
1,1-Dichloroethene	10	160 U	18 U	
1,1-Dichloroethane	10	450	18 U	
1,2-Dichloroethene (total)	10	180	18 U	
Chloroform	10	160 U	18 U	
1,2-Dichloroethane	10	160 U	18 U	
2-Butanone	10	470 J	42 J	
1,1,1-Trichloroethane	10	160	18 U	
Carbon Tetrachloride	10	160 U	18 U	
Bromodichloromethane	10	160 U	18 U	
1,2-Dichloropropane	10	160 U	18 U	
cis-1,3-Dichloropropene	10	160 U	18 U	
Trichloroethene	10	120 J	18 U	
Dibromochloromethane	10	160 U	18 U	
1,1,2-Trichloroethane	10	160 U	18 U	
Benzene	10	160 U	18 UJ	
trans-1,3-Dichloropropene	10	160 U	18 U	
Bromoform	10	160 U	18 U	
4-Methyl-2-Pentanone	10	160 U	18 U	
2-Hexanone	10	160 U	18 U	
Tetrachloroethene	10	490	3 J	
1,1,2,2-Tetrachloroethane	10	160 U	18 U	
Toluene	10	260	18 UJ	
Chlorobenzene	10	160 U	18 U	
Ethylbenzene	10	380	18 U	
Styrene	10	160 U	18 U	
Total Xylenes	10	3800	18 U	
=====				
Dilution Factor:		5.00	1.00	
Percent Solids:		32	55	
Sample Volume\Weight (ml\g):		5.00	5.00	
Associated Method Blank:		P1171.D	P1171.D	
Associated Equipment Blank:		ENQSXX1XXX94XX	ENQSXX1XXX94XX	
Associated Field Blank:		-	-	
Associated Trip Blank:		-	-	
Site: SUMP SEDIMENTS				
U: not detected J: estimated				

Table 2
Validation / Summary Table

LOCATION:	WT-101 DUP	WT-101
ISIS ID:	ENWT101XXX94XD	ENWT101XXX94XX
LAB NUMBER:	2223309	2223306
DATE SAMPLED:	10/04/94	10/04/94
DATE ANALYZED:	10/12/94	10/12/94

ANALYTE	SOW-3/90 - II	CRQL		
Chloromethane	1200	1400	U	1400 U
Bromomethane	1200	1400	U	1400 U
Vinyl Chloride	1200	1400	U	1400 U
Chloroethane	1200	1400	U	1400 U
Methylene Chloride	1200	1400	U	1400 U
Acetone	1200	1400	U	1400 U
Carbon Disulfide	1200	1400	U	1400 U
1,1-Dichloroethene	1200	1400	U	1400 U
1,1-Dichloroethane	1200	1400	U	1400 U
1,2-Dichloroethene (total)	1200	1400	U	1400 U
Chloroform	1200	1400	U	1400 U
1,2-Dichloroethane	1200	1400	U	1400 U
2-Butanone	1200	1400	U	1400 U
1,1,1-Trichloroethane	1200	1500	J	7600 J
Carbon Tetrachloride	1200	1400	U	1400 U
Bromodichloromethane	1200	1400	U	1400 U
1,2-Dichloropropane	1200	1400	U	1400 U
cis-1,3-Dichloropropene	1200	1400	U	1400 U
Trichloroethene	1200	610	J	1500 J
Dibromochloromethane	1200	1400	U	1400 U
1,1,2-Trichloroethane	1200	1400	U	1400 U
Benzene	1200	1400	U	1400 U
trans-1,3-Dichloropropene	1200	1400	U	1400 U
Bromoform	1200	1400	U	1400 U
4-Methyl-2-Pentanone	1200	1400	U	1400 U
2-Hexanone	1200	980	J	1400 U
Tetrachloroethene	1200	10000	J	43000 J
1,1,2,2-Tetrachloroethane	1200	1400	U	1400 U
Toluene	1200	1400	U	1400 U
Chlorobenzene	1200	1400	U	1400 U
Ethylbenzene	1200	1400	U	1400 U
Styrene	1200	1400	U	1400 U
Total Xylenes	1200	1400	U	520 J

Dilution Factor:	1.00	1.00
Percent Solids:	84	88
Sample Volume/Weight (ml/g):	4.00	4.00

Associated Method Blank:	N9570.D	N9570.D
Associated Equipment Blank:	ENQSXX1XXX94XX	ENQSXX1XXX94XX
Associated Field Blank:	-	-
Associated Trip Blank:	-	-

Site: WASTE-SLUDGE
U: not detected J: estimated

Table 2
Validation / Summary Table

LOCATION: WT-102 WT-103 WT-104 WT-105 WT-106						
ISIS ID: ENWT102XXX94XX ENWT103XXX94XX ENWT104XXX94XX ENWT105XXX94XX ENWT106XXX94XX						
LAB NUMBER: 2223305 2223310 2223311 2223301 2223302						
DATE SAMPLED: 10/04/94 10/04/94 10/04/94 10/05/94 10/05/94						
DATE ANALYZED: 10/12/94 10/12/94 10/12/94 10/12/94 10/12/94						
ANALYTE	SOW-3/90 - II	CRQL				
Chloromethane	10	20 U	17 U	17 U	11 U	11 U
Bromomethane	10	20 U	17 U	17 U	11 U	11 U
Vinyl Chloride	10	20 U	17 U	17 U	11 U	11 U
Chloroethane	10	20 U	17 U	17 U	11 U	11 U
Methylene Chloride	10	20 U	17 U	17 U	11 UJ	11 U
Acetone	10	20 U	190 J	17 U	6 J	11 U
Carbon Disulfide	10	20 U	4 J	17 U	2 J	11 U
1,1-Dichloroethene	10	20 U	17 U	17 U	11 U	11 U
1,1-Dichloroethane	10	26	17 U	17 U	11 U	11 U
1,2-Dichloroethene (total)	10	29	17 U	17 U	11 U	11 U
Chloroform	10	20 U	17 U	17 U	11 U	11 U
1,2-Dichloroethane	10	20 U	17 U	17 U	11 U	11 U
2-Butanone	10	20 U	14 J	17 U	5 J	11 U
1,1,1-Trichloroethane	10	75	13 J	17 U	11 U	3 J
Carbon Tetrachloride	10	9 J	17 U	17 U	11 U	11 U
Bromodichloromethane	10	20 U	17 U	17 U	11 U	11 U
1,2-Dichloropropane	10	20 U	17 U	17 U	11 U	11 U
cis-1,3-Dichloropropene	10	20 U	17 U	17 U	11 U	11 U
Trichloroethene	10	100	14 J	6 J	23 J	7 J
Dibromochloromethane	10	20 U	17 U	17 U	11 U	11 U
1,1,2-Trichloroethane	10	20 U	17 U	17 U	11 U	11 U
Benzene	10	20 U	17 U	17 U	11 U	11 U
trans-1,3-Dichloropropene	10	20 U	17 U	17 U	11 U	11 U
Bromoform	10	20 U	17 U	17 U	11 U	11 U
4-Methyl-2-Pentanone	10	20 U	17 UJ	17 UJ	11 UJ	11 UJ
2-Hexanone	10	20 U	17 UJ	17 UJ	11 UJ	11 UJ
Tetrachloroethene	10	48	10 J	37 J	43 J	31 J
1,1,2,2-Tetrachloroethane	10	20 U	17 UJ	17 UJ	11 UJ	11 UJ
Toluene	10	20 U	17 UJ	17 UJ	11 UJ	11 UJ
Chlorobenzene	10	20 U	17 UJ	17 UJ	11 UJ	11 UJ
Ethylbenzene	10	20 U	17 UJ	17 UJ	11 UJ	11 UJ
Styrene	10	20 U	17 UJ	17 UJ	11 UJ	11 UJ
Total Xylenes	10	20 U	17 UJ	17 UJ	11 UJ	11 UJ
=====						
Dilution Factor:	1.00	1.00	1.00	1.00	1.00	1.00
Percent Solids:	49	59	60	91	91	
Sample Volume/Weight (ml/g):	5.00	5.00	5.00	5.00	5.00	
Associated Method Blank:	P1148.D	P1148.D	P1148.D	P1148.D	P1148.D	
Associated Equipment Blank:	ENQSXX1XXX94XX	ENQSXX1XXX94XX	ENQSXX1XXX94XX	ENQSXX1XXX94XX	ENQSXX1XXX94XX	
Associated Field Blank:	-	-	-	-	-	
Associated Trip Blank:	-	-	-	-	-	
Site: WASTE-SLUDGE						
U: not detected J: estimated						

Table 2
Validation / Summary Table

LOCATION:	SS-101	SS-102 DUP	SS-102
ISIS ID:	ENSS101XXX94XX	ENSS102XXX94XD	ENSS102XXX94XX
LAB NUMBER:	2241101	2241105	2241102
DATE SAMPLED:	10/28/94	10/28/94	10/28/94
DATE ANALYZED:	11/03/94	11/03/94	11/03/94

ANALYTE	SOW-3/90 - II	CRQL			
Chloromethane	10		11 U	14 U	14 U
Bromomethane	10		11 U	14 U	14 U
Vinyl Chloride	10		11 U	14 U	14 U
Chloroethane	10		11 U	14 U	14 U
Methylene Chloride	10		11 U	30 U	14 U
Acetone	10		11 U	14 U	14 U
Carbon Disulfide	10		11 U	14 U	14 U
1,1-Dichloroethene	10		11 U	14 UJ	14 UJ
1,1-Dichloroethane	10		11 U	14 U	14 U
1,2-Dichloroethene (total)	10		11 U	14 U	14 U
Chloroform	10		11 U	14 U	14 U
1,2-Dichloroethane	10		11 U	14 U	14 U
2-Butanone	10		11 U	14 U	14 U
1,1,1-Trichloroethane	10		11 U	14 U	14 U
Carbon Tetrachloride	10		11 U	14 U	14 U
Bromodichloromethane	10		11 U	14 U	14 U
1,2-Dichloropropane	10		11 U	14 U	14 U
cis-1,3-Dichloropropene	10		11 U	14 U	14 U
Trichloroethene	10		11 U	14 UJ	14 UJ
Dibromochloromethane	10		11 U	14 U	14 U
1,1,2-Trichloroethane	10		11 U	14 U	14 U
Benzene	10		11 U	14 U	14 U
trans-1,3-Dichloropropene	10		11 U	14 U	14 U
Bromoform	10		11 U	14 U	14 U
4-Methyl-2-Pentanone	10		11 U	14 U	14 U
2-Hexanone	10		11 U	14 U	14 U
Tetrachloroethene	10		11 U	14 U	14 U
1,1,2,2-Tetrachloroethane	10		11 U	14 U	14 U
Toluene	10		11 U	14 U	14 U
Chlorobenzene	10		11 U	14 UJ	14 UJ
Ethylbenzene	10		11 U	14 U	14 U
Styrene	10		11 U	14 U	14 U
Total Xylenes	10		11 U	14 U	14 U

Dilution Factor:	1.00	1.00	1.00
Percent Solids:	90	74	74
Sample Volume\Weight (ml\g):	5.00	5.00	5.00
Associated Method Blank:	P1621.D	P1621.D	P1621.D
Associated Equipment Blank:	ENQS003XXX94XX	ENQS003XXX94XX	ENQS003XXX94XX
Associated Field Blank:	-	-	-
Associated Trip Blank:	-	-	-

Site: SURFACE SOILS

U: not detected J: estimated

Table 2
Validation / Summary Table

LOCATION:	BS-101	BS-102
ISIS ID:	ENBS101XX694XX	ENBS102XX694XX
LAB NUMBER:	2241106	2241107
DATE SAMPLED:	10/27/94	10/28/94
DATE ANALYZED:	11/03/94	11/03/94

ANALYTE	SOW-3/90 - II	CRQL		
Chloromethane	10	13	U	12 U
Bromomethane	10	13	U	12 U
Vinyl Chloride	10	13	U	12 U
Chloroethane	10	13	U	12 U
Methylene Chloride	10	24	U	21 U
Acetone	10	21	U	21 U
Carbon Disulfide	10	13	U	12 U
1,1-Dichloroethene	10	13	U	12 U
1,1-Dichloroethane	10	13	U	12 U
1,2-Dichloroethene (total)	10	13	U	12 U
Chloroform	10	13	U	12 U
1,2-Dichloroethane	10	13	U	12 U
2-Butanone	10	4	J	12 U
1,1,1-Trichloroethane	10	13	U	12 U
Carbon Tetrachloride	10	13	U	12 U
Bromodichloromethane	10	13	U	12 U
1,2-Dichloropropane	10	13	U	12 U
cis-1,3-Dichloropropene	10	13	U	12 U
Trichloroethene	10	13	U	12 U
Dibromochloromethane	10	13	U	12 U
1,1,2-Trichloroethane	10	13	U	12 U
Benzene	10	13	U	12 U
trans-1,3-Dichloropropene	10	13	U	12 U
Bromoform	10	13	U	12 U
4-Methyl-2-Pentanone	10	13	U	12 U
2-Hexanone	10	13	U	12 U
Tetrachloroethene	10	13	U	12 U
1,1,2,2-Tetrachloroethane	10	13	U	12 U
Toluene	10	13	U	12 U
Chlorobenzene	10	13	U	12 U
Ethylbenzene	10	13	U	12 U
Styrene	10	13	U	12 U
Total Xylenes	10	13	U	12 U

Dilution Factor:	1.00	1.00
Percent Solids:	77	81
Sample Volume\Weight (ml\g):	5.00	5.00

Associated Method Blank:	P1621.D	P1621.D
Associated Equipment Blank:	ENQS003XXX94XX	ENQS003XXX94XX
Associated Field Blank:	-	-
Associated Trip Blank:	-	-

Site: SOIL BORINGS
U: not detected J: estimated

Table 2
Validation / Summary Table

		LOCATION: MW-101 DUP		MW-101		MW-102		MW-103	
		ISIS ID: ENMW101XXX94XD		ENMW101XXX94XX		ENMW102XXX94XX		ENMW103XXX94XX	
		LAB NUMBER: 2263804		2263801		2263805		2263806	
		DATE SAMPLED: 11/30/94		11/30/94		11/30/94		11/30/94	
		DATE ANALYZED: 12/02/94		12/05/94		12/02/94		12/02/94	
ANALYTE	SOW-3/90 - II	CRQL							
Chloromethane	10		10 U	10 U	10 U	10 U	10 U	10 U	
Bromomethane	10		10 U	10 U	10 U	10 U	10 U	10 U	
Vinyl Chloride	10		10 U	10 U	10 U	10 U	10 U	10 U	
Chloroethane	10		10 U	10 U	10 U	10 U	6 J	10 U	
Methylene Chloride	10		10 UJ	10 UJ	10 U	10 U	10 U	10 U	
Acetone	10		10 U	44 UJ	11 UJ	34 UJ	10 U	10 U	
Carbon Disulfide	10		10 U	10 U	10 U	10 U	10 U	10 U	
1,1-Dichloroethene	10		10 UJ	10 UJ	10 U	10 U	10 U	10 U	
1,1-Dichloroethane	10		10 U	10 U	10 U	2 J	10 U	10 U	
1,2-Dichloroethene (total)	10		10 U	10 U	10 U	10 U	10 U	10 U	
Chloroform	10		10 U	10 U	10 U	10 U	10 U	10 U	
1,2-Dichloroethane	10		10 U	10 U	10 U	10 U	10 U	10 U	
2-Butanone	10		10 U	10 U	10 U	8 J	10 U	10 U	
1,1,1-Trichloroethane	10		10 U	10 U	10 U	10 U	10 U	10 U	
Carbon Tetrachloride	10		10 U	10 U	10 U	10 U	10 U	10 U	
Bromodichloromethane	10		10 U	10 U	10 U	10 U	10 U	10 U	
1,2-Dichloropropane	10		10 U	10 U	10 U	10 U	10 U	10 U	
cis-1,3-Dichloropropene	10		10 U	10 U	10 U	10 U	10 U	10 U	
Trichloroethene	10		10 U	10 U	10 U	10 U	10 U	10 U	
Dibromochloromethane	10		10 U	10 U	10 U	10 U	10 U	10 U	
1,1,2-Trichloroethane	10		10 U	10 U	10 U	10 U	10 U	10 U	
Benzene	10		10 U	10 U	10 U	10 U	10 U	10 U	
trans-1,3-Dichloropropene	10		10 U	10 U	10 U	10 U	10 U	10 U	
Bromoform	10		10 U	10 U	10 U	10 U	10 U	10 U	
4-Methyl-2-Pentanone	10		10 U	5 J	10 U	10 U	10 U	10 U	
2-Hexanone	10		10 U	10 U	10 U	10 U	10 U	10 U	
Tetrachloroethene	10		10 U	10 U	10 U	10 U	10 U	10 U	
1,1,2,2-Tetrachloroethane	10		10 U	10 U	10 U	10 U	10 U	10 U	
Toluene	10		10 U	10 U	10 U	10 U	10 U	10 U	
Chlorobenzene	10		10 U	10 U	10 U	10 U	10 U	10 U	
Ethylbenzene	10		10 U	10 U	10 U	10 U	10 U	10 U	
Styrene	10		10 U	10 U	10 U	10 U	10 U	10 U	
Total Xylenes	10		10 U	10 U	10 U	10 U	1 J	10 U	
=====									
Dilution Factor:		1.00	1.00	1.00	1.00	1.00	1.00	1.00	
Sample Volume\Weight (ml\g):		5.00	5.00	5.00	5.00	5.00	5.00	5.00	
Associated Method Blank:		N0496.D	N0519.D	N0496.D	N0496.D	N0496.D	N0496.D	N0496.D	
Associated Equipment Blank:		ENQS004XXX94XX	ENQS004XXX94XX	ENQS004XXX94XX	ENQS004XXX94XX	ENQS004XXX94XX	ENQS004XXX94XX	ENQS004XXX94XX	
Associated Field Blank:		-	-	-	-	-	-	-	
Associated Trip Blank:		ENQT002XXX94XX	ENQT002XXX94XX	ENQT002XXX94XX	ENQT002XXX94XX	ENQT002XXX94XX	ENQT002XXX94XX	ENQT002XXX94XX	

Site: MONITORING WELLS

U: not detected J: estimated

Table 2
Validation / Summary Table

	LOCATION:	CL-102 DUP	CL-102	CL-103	CL-104
	ISIS ID:	ENCL102XXX94XD	ENCL102XXX94XX	ENCL103XXX94XX	ENCL104XXX94XX
	LAB NUMBER:	2223317	2223314	2223318	2223312 D
	DATE SAMPLED:	10/04/94	10/04/94	10/04/94	10/05/94
	DATE EXTRACTED:	10/11/94	10/11/94	10/11/94	10/11/94
	DATE ANALYZED:	11/18/94	11/18/94	11/18/94	11/21/94
ANALYTE	SOW-3/90 - 11	CRQL			
Phenol	10	50 U	50 U	50 U	200 U
bis(2-Chloroethyl)ether	10	50 U	50 U	50 U	200 U
2-Chlorophenol	10	50 U	50 U	50 U	200 U
1,3-Dichlorobenzene	10	50 U	50 U	50 U	200 U
1,4-Dichlorobenzene	10	50 U	50 U	50 U	200 U
1,2-Dichlorobenzene	10	50 U	50 U	50 U	200 U
2-Methylphenol	10	50 U	50 U	50 U	200 U
2,2'-oxybis(1-Chloropropane)	10	50 U	50 U	50 U	200 U
4-Methylphenol	10	50 U	50 U	50 U	31 J
N-Nitroso-di-n-propylamine	10	50 U	50 U	50 U	200 U
Hexachloroethane	10	50 U	50 U	50 U	200 U
Nitrobenzene	10	50 U	50 U	50 U	200 U
Isophorone	10	50 U	50 U	50 U	200 U
2-Nitrophenol	10	50 U	50 U	50 U	200 U
2,4-Dimethylphenol	10	50 U	50 U	50 U	200 U
bis(2-Chloroethoxy)methane	10	50 U	50 U	50 U	200 U
2,4-Dichlorophenol	10	50 U	50 U	50 U	200 U
1,2,4-Trichlorobenzene	10	50 U	50 U	50 U	200 U
Naphthalene	10	50 U	50 U	50 U	200 U
4-Chloroaniline	10	50 U	50 U	50 U	200 U
Hexachlorobutadiene	10	50 U	50 U	50 U	200 U
4-Chloro-3-Methylphenol	10	50 U	50 U	50 U	200 U
2-Methylnaphthalene	10	50 U	50 U	50 U	22 J
Hexachlorocyclopentadiene	10	50 U	50 U	50 U	200 U
2,4,6-Trichlorophenol	10	50 U	50 U	50 U	200 U
2,4,5-Trichlorophenol	25	120 U	120 U	120 U	500 U
2-Chloronaphthalene	10	50 U	50 U	50 U	200 U
2-Nitroaniline	25	120 U	120 U	120 U	500 U
Dimethylphthalate	10	50 U	50 U	50 U	200 U
Acenaphthylene	10	50 U	50 U	50 U	200 U
2,6-Dinitrotoluene	10	50 U	50 U	50 U	200 U

Site: SUMP LIQUIDS

U: not detected J: estimated

Table 2
Validation / Summary Table

LOCATION:	CL-102 DUP	CL-102	CL-103	CL-104
ISIS ID:	ENCL102XXX94XD	ENCL102XXX94XX	ENCL103XXX94XX	ENCL104XXX94XX
LAB NUMBER:	2223317	2223314	2223318	2223312 D
DATE SAMPLED:	10/04/94	10/04/94	10/04/94	10/05/94
DATE EXTRACTED:	10/11/94	10/11/94	10/11/94	10/11/94
DATE ANALYZED:	11/18/94	11/18/94	11/18/94	11/21/94

ANALYTE	SOW-3/90 - II	CRQL				
3-Nitroaniline	25	120 U	120 U	120 U	500 U	
Acenaphthene	10	50 U	50 U	50 U	200 U	
2,4-Dinitrophenol	25	120 UJ	120 UJ	120 UJ		R
4-Nitrophenol	25	120 U	120 U	120 U	500 U	
Dibenzofuran	10	50 U	50 U	50 U	200 U	
2,4-Dinitrotoluene	10	50 U	50 U	50 U	200 U	
Diethylphthalate	10	50 U	50 U	50 U	200 U	
4-Chlorophenyl-phenylether	10	50 U	50 U	50 U	200 U	
Fluorene	10	50 U	50 U	50 U	200 U	
4-Nitroaniline	25	120 UJ	120 UJ	120 UJ	500 U	
4,6-Dinitro-2-methylphenol	25	120 U	120 U	120 U	500 UJ	
N-Nitrosodiphenylamine	10	50 U	50 U	50 U	200 U	
4-Bromophenyl-phenylether	10	50 U	50 U	50 U	200 U	
Hexachlorobenzene	10	50 U	50 U	50 U	200 U	
Pentachlorophenol	25	120 U	120 U	120 U	180 J	
Phenanthrene	10	50 U	50 U	50 U	200 U	
Anthracene	10	50 U	50 U	50 U	200 U	
Carbazole	10	50 U	50 U	50 U	200 U	
Di-n-butylphthalate	10	50 U	50 U	50 U	200 U	
Fluoranthene	10	50 U	50 U	50 U	200 U	
Pyrene	10	6 J	50 U	50 U	200 U	
Butylbenzylphthalate	10	38 J	38 J	50 U	30 J	
3,3'-Dichlorobenzidine	10	50 U	50 U	50 U	200 UJ	
Benzo(a)Anthracene	10	50 U	50 U	50 U	200 U	
Chrysene	10	50 U	50 U	50 U	200 U	
bis(2-Ethylhexyl)phthalate	10	380	290	50 U	400 U	
Di-n-octylphthalate	10	130 J	40 J	50 UJ	63 J	
Benzo(b)Fluoranthene	10	50 U	50 U	50 UJ	200 UJ	
Benzo(k)Fluoranthene	10	50 U	50 U	50 UJ	200 UJ	
Benzo(a)Pyrene	10	50 U	50 U	50 UJ	200 UJ	
Indeno(1,2,3-c,d)Pyrene	10	50 U	50 U	50 UJ	200 UJ	
Dibenz(a,h)Anthracene	10	50 U	50 U	50 UJ	200 UJ	
Benzo(g,h,i)perylene	10	50 U	50 U	50 UJ	200 UJ	
=====						
Dilution Factor:	5.00	5.00	5.00	20.0		
Sample Volume\Weight (ml\g):	1000	1000	1000	1000		
Associated Method Blank:	S1523.D	S1523.D	S1523.D	S1523.D		
Associated Equipment Blank:	-	-	-	-		
Associated Field Blank:	-	-	-	-		

Site: SUMP LIQUIDS

U: not detected J: estimated

Table 2
Validation / Summary Table

LOCATION:	CD-101	CD-103
ISIS ID:	ENCD101XXX94XX	ENCD103XXX94XX
LAB NUMBER:	2223303	2223304
DATE SAMPLED:	10/04/94	10/04/94
DATE EXTRACTED:	10/08/94	10/08/94
DATE ANALYZED:	11/17/94	11/17/94

ANALYTE	SOW-3/90 - II	CRQL			
Phenol	330	21000	UJ	30000	U
bis(2-Chloroethyl)ether	330	21000	UJ	30000	U
2-Chlorophenol	330	21000	UJ	30000	U
1,3-Dichlorobenzene	330	21000	UJ	30000	U
1,4-Dichlorobenzene	330	21000	UJ	30000	U
1,2-Dichlorobenzene	330	21000	UJ	30000	U
2-Methylphenol	330	21000	UJ	30000	U
2,2'-oxybis(1-Chloropropane)	330	21000	UJ	30000	U
4-Methylphenol	330	21000	UJ	30000	U
N-Nitroso-di-n-propylamine	330	21000	UJ	30000	U
Hexachloroethane	330	21000	UJ	30000	U
Nitrobenzene	330	21000	UJ	30000	U
Isophorone	330	21000	UJ	30000	U
2-Nitrophenol	330	21000	UJ	30000	U
2,4-Dimethylphenol	330	21000	UJ	30000	U
bis(2-Chloroethoxy)methane	330	21000	UJ	30000	U
2,4-Dichlorophenol	330	21000	UJ	30000	U
1,2,4-Trichlorobenzene	330	21000	UJ	30000	U
Naphthalene	330	21000	UJ	30000	U
4-Chloroaniline	330	21000	UJ	30000	U
Hexachlorobutadiene	330	21000	UJ	30000	U
4-Chloro-3-Methylphenol	330	21000	UJ	30000	U
2-Methylnaphthalene	330	21000	UJ	30000	U
Hexachlorocyclopentadiene	330	21000	UJ	30000	U
2,4,6-Trichlorophenol	330	21000	UJ	30000	U
2,4,5-Trichlorophenol	800	50000	UJ	73000	U
2-Chloronaphthalene	330	21000	UJ	30000	U
2-Nitroaniline	800	50000	UJ	73000	U
Dimethylphthalate	330	21000	UJ	30000	U
Acenaphthylene	330	21000	UJ	30000	U
2,6-Dinitrotoluene	330	21000	UJ	30000	U

Site: SUMP SEDIMENTS

U: not detected J: estimated

Table 2
Validation / Summary Table

LOCATION:	CD-101	CD-103
ISIS ID:	ENCD101XXX94XX	ENCD103XXX94XX
LAB NUMBER:	2223303	2223304
DATE SAMPLED:	10/04/94	10/04/94
DATE EXTRACTED:	10/08/94	10/08/94
DATE ANALYZED:	11/17/94	11/17/94

ANALYTE	SOW-3/90 - II	CRQL				
3-Nitroaniline	800	50000	UJ	73000	U	
Acenaphthene	330	21000	UJ	30000	U	
2,4-Dinitrophenol	800	50000	UJ	73000	UJ	
4-Nitrophenol	800	50000	UJ	73000	U	
Dibenzofuran	330	21000	UJ	30000	U	
2,4-Dinitrotoluene	330	21000	UJ	30000	U	
Diethylphthalate	330	21000	UJ	30000	U	
4-Chlorophenyl-phenylether	330	21000	UJ	30000	U	
Fluorene	330	21000	UJ	30000	U	
4-Nitroaniline	800	50000	UJ	73000	U	
4,6-Dinitro-2-methylphenol	800	50000	UJ	73000	UJ	
N-Nitrosodiphenylamine	330	21000	UJ	30000	U	
4-Bromophenyl-phenylether	330	21000	UJ	30000	U	
Hexachlorobenzene	330	21000	UJ	30000	U	
Pentachlorophenol	800	50000	UJ	73000	U	
Phenanthrene	330	21000	UJ	30000	U	
Anthracene	330	21000	UJ	30000	U	
Carbazole	330	21000	UJ	30000	U	
Di-n-butylphthalate	330	21000	UJ	30000	U	
Fluoranthene	330	21000	UJ	30000	U	
Pyrene	330	21000	UJ	30000	U	
Butylbenzylphthalate	330	3000	J	30000	U	
3,3'-Dichlorobenzidine	330	21000	UJ	30000	U	
Benzo(a)Anthracene	330	21000	UJ	30000	U	
Chrysene	330	21000	UJ	30000	U	
bis(2-Ethylhexyl)phthalate	330	46000	UJ	30000	U	
Di-n-octylphthalate	330	21000	UJ	30000	U	
Benzo(b)Fluoranthene	330	21000	UJ	30000	U	
Benzo(k)Fluoranthene	330	21000	UJ	30000	U	
Benzo(a)Pyrene	330	21000	UJ	30000	U	
Indeno(1,2,3-c,d)Pyrene	330	21000	UJ	30000	U	
Dibenz(a,h)Anthracene	330	21000	UJ	30000	U	
Benzo(g,h,i)perylene	330	21000	UJ	30000	U	

Dilution Factor:	20.0	50.0
Percent Solids:	32	55
Sample Volume/Weight (ml/g):	30.0	30.0

Associated Method Blank:	S1503.D	S1503.D
Associated Equipment Blank:	ENQSXX1XXX94XX	ENQSXX1XXX94XX
Associated Field Blank:	-	-

Site: SUMP SEDIMENTS
U: not detected J: estimated

Table 2
Validation / Summary Table

			LOCATION:	WT-101 DUP	WT-101	WT-102	WT-103	WT-104	WT-105	WT-106
			ISIS ID:	ENWT101XXX94XD	ENWT101XXX94XX	ENWT102XXX94XX	ENWT103XXX94XX	ENWT104XXX94XX	ENWT105XXX94XX	ENWT106XXX94XX
			LAB NUMBER:	2223309	2223306	2223305	2223310	2223311	2223301	2223302
			DATE SAMPLED:	10/04/94	10/04/94	10/04/94	10/04/94	10/04/94	10/05/94	10/05/94
			DATE EXTRACTED:	10/08/94	10/08/94	10/08/94	10/08/94	10/08/94	10/08/94	10/08/94
			DATE ANALYZED:	11/17/94	11/17/94	11/17/94	11/17/94	11/17/94	11/17/94	11/17/94
ANALYTE	SOW-3/90 - II	CRQL								
Phenol	330	R				69000 UJ	11000 U	11000 U	7300 U	37000 U
bis(2-Chloroethyl)ether	330	7900 U			150000 U	69000 UJ	11000 U	11000 U	7300 U	37000 U
2-Chlorophenol	330	R				69000 UJ	11000 U	11000 U	7300 U	37000 U
1,3-Dichlorobenzene	330	7900 U			150000 U	69000 UJ	11000 U	11000 U	7300 U	37000 U
1,4-Dichlorobenzene	330	R				69000 UJ	11000 U	11000 U	7300 U	37000 U
1,2-Dichlorobenzene	330	7900 U			150000 U	69000 UJ	11000 U	11000 U	7300 U	37000 U
2-Methylphenol	330	7900 U			150000 U	69000 UJ	11000 U	11000 U	7300 U	37000 U
2,2'-oxybis(1-Chloropropane)	330	7900 U			150000 U	69000 UJ	11000 U	11000 U	7300 U	37000 U
4-Methylphenol	330	7900 U			150000 U	69000 UJ	11000 U	11000 U	7300 U	37000 U
N-Nitroso-di-n-propylamine	330	R				69000 UJ	11000 U	11000 U	7300 U	37000 U
Hexachloroethane	330	7900 U			150000 U	69000 UJ	11000 U	11000 U	7300 U	37000 U
Nitrobenzene	330	7900 U			150000 U	69000 UJ	11000 U	11000 U	7300 U	37000 U
Isophorone	330	7900 U			150000 U	69000 UJ	11000 U	11000 U	7300 U	37000 U
2-Nitrophenol	330	7900 U			150000 U	69000 UJ	11000 U	11000 U	7300 U	37000 U
2,4-Dimethylphenol	330	7900 U			150000 U	69000 UJ	11000 U	11000 U	7300 U	37000 U
bis(2-Chloroethoxy)methane	330	7900 U			150000 U	69000 UJ	11000 U	11000 U	7300 U	37000 U
2,4-Dichlorophenol	330	7900 U			150000 U	69000 UJ	11000 U	11000 U	7300 U	37000 U
1,2,4-Trichlorobenzene	330	R				69000 UJ	11000 U	11000 U	7300 U	37000 U
Naphthalene	330	7900 U			150000 U	69000 UJ	11000 U	11000 U	7300 U	37000 U
4-Chloroaniline	330	7900 U			150000 U	69000 UJ	11000 U	11000 U	7300 U	37000 U
Hexachlorobutadiene	330	7900 U			150000 U	69000 UJ	11000 U	11000 U	7300 U	37000 U
4-Chloro-3-Methylphenol	330	R				69000 UJ	11000 U	11000 U	7300 U	37000 U
2-Methylnaphthalene	330	7900 U			150000 U	69000 UJ	11000 U	11000 U	7300 U	37000 U
Hexachlorocyclopentadiene	330	7900 U			150000 U	69000 UJ	11000 U	11000 U	7300 U	37000 U
2,4,6-Trichlorophenol	330	7900 U			150000 U	69000 UJ	11000 U	11000 U	7300 U	37000 U
2,4,5-Trichlorophenol	800	19000 U			360000 U	170000 UJ	27000 U	27000 U	18000 U	88000 U
2-Chloronaphthalene	330	7900 U			150000 U	69000 UJ	11000 U	11000 U	7300 U	37000 U
2-Nitroaniline	800	19000 U			360000 U	170000 UJ	27000 U	27000 U	18000 U	88000 U
Dimethylphthalate	330	7900 U			150000 U	69000 UJ	11000 U	11000 U	7300 U	37000 U
Acenaphthylene	330	7900 U			150000 U	69000 UJ	11000 U	11000 U	7300 U	37000 U
2,6-Dinitrotoluene	330	7900 U			150000 U	69000 UJ	11000 U	11000 U	7300 U	37000 U

Site: WASTE-SLUDGE

U: not detected R: unusable J: estimated

Table 2
Validation / Summary Table

LOCATION:		WT-101 DUP		WT-101		WT-102		WT-103		WT-104		WT-105		WT-106	
ISIS ID:		ENWT101XXX94XD		ENWT101XXX94XX		ENWT102XXX94XX		ENWT103XXX94XX		ENWT104XXX94XX		ENWT105XXX94XX		ENWT106XXX94XX	
LAB NUMBER:		2223309		2223306		2223305		2223310		2223311		2223301		2223302	
DATE SAMPLED:		10/04/94		10/04/94		10/04/94		10/04/94		10/04/94		10/05/94		10/05/94	
DATE EXTRACTED:		10/08/94		10/08/94		10/08/94		10/08/94		10/08/94		10/08/94		10/08/94	
DATE ANALYZED:		11/17/94		11/17/94		11/17/94		11/17/94		11/17/94		11/17/94		11/17/94	
ANALYTE	SOW-3/90 - II	CRQL													
3-Nitroaniline	800	19000	U	360000	U	170000	UJ	27000	U	27000	U	18000	U	88000	U
Acenaphthene	330		R		R	69000	UJ	11000	U	11000	U	7300	U	37000	U
2,4-Dinitrophenol	800	19000	UJ	360000	UJ	170000	UJ	27000	UJ	27000	UJ	18000	UJ	88000	UJ
4-Nitrophenol	800		R		R	170000	UJ	27000	U	27000	U	18000	U	88000	U
Dibenzofuran	330	7900	U	150000	U	69000	UJ	11000	U	11000	U	7300	U	37000	U
2,4-Dinitrotoluene	330		R		R	69000	UJ	11000	U	11000	U	7300	U	37000	U
Diethylphthalate	330	7900	U	150000	U	69000	UJ	11000	U	11000	U	7300	U	37000	U
4-Chlorophenyl-phenylether	330	7900	U	150000	U	69000	UJ	11000	U	11000	U	7300	U	37000	U
Fluorene	330	7900	U	150000	U	69000	UJ	11000	U	11000	U	7300	U	37000	U
4-Nitroaniline	800	19000	U	360000	U	170000	UJ	27000	U	27000	U	18000	U	88000	U
4,6-Dinitro-2-methylphenol	800	19000	UJ	360000	UJ	170000	UJ	27000	UJ	27000	UJ	18000	UJ	88000	UJ
N-Nitrosodiphenylamine	330	7900	U	150000	U	69000	UJ	11000	U	11000	U	7300	U	37000	U
4-Bromophenyl-phenylether	330	7900	U	150000	U	69000	UJ	11000	U	11000	U	7300	U	37000	U
Hexachlorobenzene	330	7900	U	150000	U	69000	UJ	11000	U	11000	U	7300	U	37000	U
Pentachlorophenol	800		R		R	170000	UJ	27000	U	27000	U	18000	U	7400	J
Phenanthrene	330	1100	J	150000	U	69000	UJ	11000	U	11000	U	7300	U	37000	U
Anthracene	330	7900	U	150000	U	69000	UJ	11000	U	11000	U	7300	U	37000	U
Carbazole	330	7900	U	150000	U	69000	UJ	11000	U	11000	U	7300	U	37000	U
Di-n-butylphthalate	330	7900	U	150000	U	69000	UJ	11000	U	11000	U	7300	U	37000	U
Fluoranthene	330	940	J	150000	U	69000	UJ	11000	U	11000	U	7300	U	37000	U
Pyrene	330	1800	J		R	69000	UJ	11000	U	11000	U	7300	U	37000	U
Butylbenzylphthalate	330	7900	U	150000	U	69000	UJ	11000	U	11000	U	7300	U	37000	U
3,3'-Dichlorobenzidine	330	7900	U	150000	U	69000	UJ	11000	U	11000	U	7300	U	37000	U
Benzo(a)Anthracene	330	7900	U	150000	U	69000	UJ	11000	U	11000	U	7300	U	37000	U
Chrysene	330	940	J	150000	U	69000	UJ	11000	U	11000	U	7300	U	37000	U
bis(2-Ethylhexyl)phthalate	330	7900	U	150000	U	69000	UJ	11000	U	11000	U	7300	U	37000	U
Di-n-octylphthalate	330	7900	U	150000	U	69000	UJ	11000	U	11000	U	7300	U	37000	U
Benzo(b)Fluoranthene	330	7900	U	150000	U	69000	UJ	11000	U	11000	U	7300	U	37000	U
Benzo(k)Fluoranthene	330	7900	U	150000	U	69000	UJ	11000	U	11000	U	7300	U	37000	U
Benzo(a)Pyrene	330	7900	U	150000	U	69000	UJ	11000	U	11000	U	7300	U	37000	U
Indeno(1,2,3-c,d)Pyrene	330	7900	U	150000	U	69000	UJ	11000	U	11000	U	7300	U	37000	U
Dibenz(a,h)Anthracene	330	7900	U	150000	U	69000	UJ	11000	U	11000	U	7300	U	37000	U
Benzo(g,h,i)perylene	330	7900	U	150000	U	69000	UJ	11000	U	11000	U	7300	U	37000	U
=====															
Dilution Factor:		20.0		400		100		20.0		20.0		20.0		100	
Percent Solids:		84		88		48		59		60		91		91	
Sample Volume\Weight (ml\g):		30.0		30.0		30.0		30.0		30.0		30.0		30.0	
Associated Method Blank:		S1503.D		S1503.D		S1503.D		S1503.D		S1503.D		S1503.D		S1503.D	
Associated Equipment Blank:		ENQSXX1XXX94XX		ENQSXX1XXX94XX		ENQSXX1XXX94XX		ENQSXX1XXX94XX		ENQSXX1XXX94XX		ENQSXX1XXX94XX		ENQSXX1XXX94XX	
Associated Field Blank:		-		-		-		-		-		-		-	

Site: WASTE-SLUDGE

U: not detected R: unusable J: estimated

Table 2
Validation / Summary Table

LOCATION:	SS-101	SS-102 DUP	SS-102
ISIS ID:	ENSS101XXX94XX	ENSS102XXX94XD	ENSS102XXX94XX
LAB NUMBER:	2241101	2241105	2241102
DATE SAMPLED:	10/28/94	10/28/94	10/28/94
DATE EXTRACTED:	11/01/94	11/01/94	11/01/94
DATE ANALYZED:	12/03/94	12/03/94	12/03/94

ANALYTE	SOW-3/90 - 11	CRQL			
Phenol	330	1800 U	900 U	900 U	
bis(2-Chloroethyl)ether	330	1800 U	900 U	900 U	
2-Chlorophenol	330	1800 U	900 U	900 U	
1,3-Dichlorobenzene	330	1800 U	900 U	900 U	
1,4-Dichlorobenzene	330	1800 U	900 UJ	900 UJ	
1,2-Dichlorobenzene	330	1800 U	900 U	900 U	
2-Methylphenol	330	1800 U	900 U	900 U	
2,2'-oxybis(1-Chloropropane)	330	1800 U	900 U	900 U	
4-Methylphenol	330	1800 U	900 U	900 U	
N-Nitroso-di-n-propylamine	330	1800 U	900 U	900 U	
Hexachloroethane	330	1800 U	900 U	900 U	
Nitrobenzene	330	1800 U	900 U	900 U	
Isophorone	330	1800 U	900 U	900 U	
2-Nitrophenol	330	1800 U	900 U	900 U	
2,4-Dimethylphenol	330	1800 U	900 U	900 U	
bis(2-Chloroethoxy)methane	330	1800 U	900 U	900 U	
2,4-Dichlorophenol	330	1800 U	900 U	900 U	
1,2,4-Trichlorobenzene	330	1800 U	900 U	900 U	
Naphthalene	330	1800 U	900 U	900 U	
4-Chloroaniline	330	1800 U	900 U	900 U	
Hexachlorobutadiene	330	1800 U	900 U	900 U	
4-Chloro-3-Methylphenol	330	1800 U	900 U	900 U	
2-Methylnaphthalene	330	1800 U	900 U	900 U	
Hexachlorocyclopentadiene	330	R	R	R	
2,4,6-Trichlorophenol	330	1800 U	900 U	900 U	
2,4,5-Trichlorophenol	800	4400 U	2200 U	2200 U	
2-Chloronaphthalene	330	1800 U	900 U	900 U	
2-Nitroaniline	800	4400 U	2200 U	2200 U	
Dimethylphthalate	330	1800 U	900 U	900 U	
Acenaphthylene	330	1800 U	900 U	900 U	
2,6-Dinitrotoluene	330	1800 U	900 U	900 U	

Site: SURFACE SOILS

U: not detected J: estimated R: unusable

Table 2
Validation / Summary Table

LOCATION:	SS-101	SS-102 DUP	SS-102
ISIS ID:	ENSS101XXX94XX	ENSS102XXX94XD	ENSS102XXX94XX
LAB NUMBER:	2241101	2241105	2241102
DATE SAMPLED:	10/28/94	10/28/94	10/28/94
DATE EXTRACTED:	11/01/94	11/01/94	11/01/94
DATE ANALYZED:	12/03/94	12/03/94	12/03/94

ANALYTE	SOW-3/90 - II	CRQL			
3-Nitroaniline	800	4400 U	2200 U	2200 U	
Acenaphthene	330	470 J	900 U	900 U	
2,4-Dinitrophenol	800	4400 U	2200 U	2200 U	
4-Nitrophenol	800	4400 U	2200 U	2200 U	
Dibenzofuran	330	290 J	900 U	900 U	
2,4-Dinitrotoluene	330	1800 U	900 U	900 U	
Diethylphthalate	330	1800 U	900 U	900 U	
4-Chlorophenyl-phenylether	330	1800 U	900 U	900 U	
Fluorene	330	560 J	900 U	900 U	
4-Nitroaniline	800	4400 U	2200 U	2200 U	
4,6-Dinitro-2-methylphenol	800	4400 U	2200 U	2200 U	
N-Nitrosodiphenylamine	330	1800 U	900 U	900 U	
4-Bromophenyl-phenylether	330	1800 U	900 U	900 U	
Hexachlorobenzene	330	1800 U	900 U	900 U	
Pentachlorophenol	800	4400 U	R	R	
Phenanthrene	330	6100	220 J	450 J	
Anthracene	330	1100 J	900 U	900 U	
Carbazole	330	640 J	900 U	900 U	
Di-n-butylphthalate	330	1800 U	900 U	900 U	
Fluoranthene	330	9400	320 J	680 J	
Pyrene	330	7900	300 J	620 J	
Butylbenzylphthalate	330	1800 U	900 U	900 U	
3,3'-Dichlorobenzidine	330	1800 U	900 U	900 U	
Benzo(a)Anthracene	330	3900	160 J	350 J	
Chrysene	330	4400	200 J	450 J	
bis(2-Ethylhexyl)phthalate	330	1800 U	900 U	900 U	
Di-n-octylphthalate	330	1800 U	900 U	900 U	
Benzo(b)Fluoranthene	330	3200	130 J	270 J	
Benzo(k)Fluoranthene	330	2100	900 U	170 J	
Benzo(a)Pyrene	330	2400	95 J	210 J	
Indeno(1,2,3-c,d)Pyrene	330	720 J	900 U	120 J	
Dibenz(a,h)Anthracene	330	1800 U	900 U	900 U	
Benzo(g,h,i)perylene	330	600 J	900 U	120 J	

Dilution Factor:	5.00	2.00	2.00
Percent Solids:	90	74	74
Sample Volume\Weight (ml\g):	30.0	30.0	30.0
Associated Method Blank:	R1715.D	R1715.D	R1715.D
Associated Equipment Blank:	ENQS003XXX94XX	ENQS003XXX94XX	ENQS003XXX94XX
Associated Field Blank:	-	-	-

Site: SURFACE SOILS

U: not detected J: estimated R: unusable

Table 2
Validation / Summary Table

LOCATION:	BS-101	BS-102
ISIS ID:	ENBS101XX694XX	ENBS102XX694XX
LAB NUMBER:	2241106	2241107
DATE SAMPLED:	10/27/94	10/28/94
DATE EXTRACTED:	11/01/94	11/01/94
DATE ANALYZED:	12/03/94	12/03/94

ANALYTE	SOW-3/90 - 11	CRQL			
Phenol	330	860	U	410	U
bis(2-Chloroethyl)ether	330	860	U	410	U
2-Chlorophenol	330	860	U	410	U
1,3-Dichlorobenzene	330	860	U	410	U
1,4-Dichlorobenzene	330	860	U	410	U
1,2-Dichlorobenzene	330	860	U	410	U
2-Methylphenol	330	860	U	410	U
2,2'-oxybis(1-Chloropropane)	330	860	U	410	U
4-Methylphenol	330	860	U	410	U
N-Nitroso-di-n-propylamine	330	860	U	410	U
Hexachloroethane	330	860	U	410	U
Nitrobenzene	330	860	U	410	U
Isophorone	330	860	U	410	U
2-Nitrophenol	330	860	U	410	U
2,4-Dimethylphenol	330	860	U	410	U
bis(2-Chloroethoxy)methane	330	860	U	410	U
2,4-Dichlorophenol	330	860	U	410	U
1,2,4-Trichlorobenzene	330	860	U	410	U
Naphthalene	330	860	U	410	U
4-Chloroaniline	330	860	U	410	U
Hexachlorobutadiene	330	860	U	410	U
4-Chloro-3-Methylphenol	330	860	U	410	U
2-Methylnaphthalene	330	860	U	410	U
Hexachlorocyclopentadiene	330		R		R
2,4,6-Trichlorophenol	330	860	U	410	U
2,4,5-Trichlorophenol	800	2100	U	990	U
2-Chloronaphthalene	330	860	U	410	U
2-Nitroaniline	800	2100	U	990	U
Dimethylphthalate	330	860	U	410	U
Acenaphthylene	330	860	U	410	U
2,6-Dinitrotoluene	330	860	U	410	U

Site: SOIL BORINGS

U: not detected R: unusable J: estimated

Table 2
Validation / Summary Table

LOCATION:	BS-101	BS-102
ISIS ID:	ENBS101XX694XX	ENBS102XX694XX
LAB NUMBER:	2241106	2241107
DATE SAMPLED:	10/27/94	10/28/94
DATE EXTRACTED:	11/01/94	11/01/94
DATE ANALYZED:	12/03/94	12/03/94

ANALYTE	SOW-3/90 - II	CRQL		
3-Nitroaniline	800	2100	U	990 U
Acenaphthene	330	860	U	410 U
2,4-Dinitrophenol	800	2100	U	990 U
4-Nitrophenol	800	2100	U	990 U
Dibenzofuran	330	860	U	410 U
2,4-Dinitrotoluene	330	860	U	410 U
Diethylphthalate	330	860	U	410 U
4-Chlorophenyl-phenylether	330	860	U	410 U
Fluorene	330	860	U	410 U
4-Nitroaniline	800	2100	U	990 U
4,6-Dinitro-2-methylphenol	800	2100	U	990 U
N-Nitrosodiphenylamine	330	860	U	410 U
4-Bromophenyl-phenylether	330	860	U	410 U
Hexachlorobenzene	330	860	U	410 U
Pentachlorophenol	800	2100	U	990 U
Phenanthrene	330	300	J	110 J
Anthracene	330	860	U	410 U
Carbazole	330	860	U	410 U
Di-n-butylphthalate	330	860	U	410 U
Fluoranthene	330	260	J	200 J
Pyrene	330	240	J	180 J
Butylbenzylphthalate	330	860	U	410 U
3,3'-Dichlorobenzidine	330	860	U	410 U
Benzo(a)Anthracene	330	130	J	100 J
Chrysene	330	140	J	120 J
bis(2-Ethylhexyl)phthalate	330	860	U	410 U
Di-n-octylphthalate	330	860	U	410 U
Benzo(b)Fluoranthene	330	860	U	85 J
Benzo(k)Fluoranthene	330	860	U	61 J
Benzo(a)Pyrene	330	860	U	65 J
Indeno(1,2,3-c,d)Pyrene	330	860	U	44 J
Dibenz(a,h)Anthracene	330	860	U	410 U
Benzo(g,h,i)perylene	330	860	U	45 J

Dilution Factor:	2.00	1.00
Percent Solids:	77	81
Sample Volume\Weight (ml\g):	30.0	30.0

Associated Method Blank:	R1715.D	R1715.D
Associated Equipment Blank:	ENQS003XXX94XX	ENQS003XXX94XX
Associated Field Blank:	-	-

Site: SOIL BORINGS

U: not detected R: unusable J: estimated

Table 2
Validation / Summary Table

LOCATION:	MW-101 DUP	MW-101	MW-102	MW-103
ISIS ID:	ENMW101XXX94XD	ENMW101XXX94XX	ENMW102XXX94XX	ENMW103XXX94XX
LAB NUMBER:	2263804	2263801	2263805	2263806
DATE SAMPLED:	11/30/94	11/30/94	11/30/94	11/30/94
DATE EXTRACTED:	12/06/94	12/06/94	12/06/94	12/06/94
DATE ANALYZED:	12/22/94	12/22/94	12/22/94	12/22/94

ANALYTE	SOW-3/90 - II	CRQL				
Phenol	10	10 U	10 U	10 U	10 U	10 U
bis(2-Chloroethyl)ether	10	10 U	10 U	10 U	10 U	10 U
2-Chlorophenol	10	10 U	10 U	10 U	10 U	10 U
1,3-Dichlorobenzene	10	10 U	10 U	10 U	10 U	10 U
1,4-Dichlorobenzene	10	10 U	10 U	10 U	10 U	10 U
1,2-Dichlorobenzene	10	10 U	10 U	10 U	10 U	10 U
2-Methylphenol	10	10 U	10 U	10 U	2 J	10 U
2,2'-oxybis(1-Chloropropane)	10	10 U	10 U	10 U	10 U	10 U
4-Methylphenol	10	10 U	10 U	10 U	15	10 U
N-Nitroso-di-n-propylamine	10	10 U	10 U	10 U	10 U	10 U
Hexachloroethane	10	10 U	10 U	10 U	10 U	10 U
Nitrobenzene	10	10 U	10 U	10 U	10 U	10 U
Isophorone	10	10 U	10 U	10 U	10 U	10 U
2-Nitrophenol	10	10 U	10 U	10 U	10 U	10 U
2,4-Dimethylphenol	10	10 U	10 U	10 U	1 J	10 U
bis(2-Chloroethoxy)methane	10	10 U	10 U	10 U	10 U	10 U
2,4-Dichlorophenol	10	10 U	10 U	10 U	10 U	10 U
1,2,4-Trichlorobenzene	10	10 U	10 U	10 U	10 U	10 U
Naphthalene	10	10 U	10 U	10 U	14	10 U
4-Chloroaniline	10	10 U	10 U	10 U	10 U	10 U
Hexachlorobutadiene	10	10 U	10 U	10 U	10 U	10 U
4-Chloro-3-Methylphenol	10	10 U	10 U	10 U	10 U	10 U
2-Methylnaphthalene	10	10 U	10 U	10 U	4 J	10 U
Hexachlorocyclopentadiene	10	10 U	10 U	10 U	10 U	10 U
2,4,6-Trichlorophenol	10	10 U	10 U	10 U	10 U	10 U
2,4,5-Trichlorophenol	25	25 U	25 U	25 U	25 U	25 U
2-Chloronaphthalene	10	10 U	10 U	10 U	10 U	10 U
2-Nitroaniline	25	25 U	25 U	25 U	25 U	25 U
Dimethylphthalate	10	10 U	10 U	10 U	10 U	10 U
Acenaphthylene	10	10 U	10 U	10 U	10 U	10 U
2,6-Dinitrotoluene	10	10 U	10 U	10 U	10 U	10 U

Site: MONITORING WELLS
U: not detected J: estimated

Table 2
Validation / Summary Table

LOCATION:	MW-101 DUP	MW-101	MW-102	MW-103
ISIS ID:	ENMW101XXX94XD	ENMW101XXX94XX	ENMW102XXX94XX	ENMW103XXX94XX
LAB NUMBER:	2263804	2263801	2263805	2263806
DATE SAMPLED:	11/30/94	11/30/94	11/30/94	11/30/94
DATE EXTRACTED:	12/06/94	12/06/94	12/06/94	12/06/94
DATE ANALYZED:	12/22/94	12/22/94	12/22/94	12/22/94

ANALYTE	SOW-3/90 - II	CRQL				
3-Nitroaniline	25	25 UJ	25 UJ	25 UJ	25 UJ	25 UJ
Acenaphthene	10	10 U	10 U	10 U	10 U	10 U
2,4-Dinitrophenol	25	25 UJ	25 UJ	25 UJ	25 UJ	25 UJ
4-Nitrophenol	25	25 U	25 U	25 U	25 U	25 U
Dibenzofuran	10	10 U	10 U	10 U	10 U	1 J
2,4-Dinitrotoluene	10	10 U	10 U	10 U	10 U	10 U
Diethylphthalate	10	10 U	10 U	10 U	10 U	10 U
4-Chlorophenyl-phenylether	10	10 U	10 U	10 U	10 U	10 U
Fluorene	10	10 U	10 U	10 U	10 U	2 J
4-Nitroaniline	25	25 U	25 U	25 U	25 U	25 U
4,6-Dinitro-2-methylphenol	25	25 UJ	25 UJ	25 UJ	25 UJ	25 UJ
N-Nitrosodiphenylamine	10	10 UJ	10 UJ	10 UJ	10 UJ	29 J
4-Bromophenyl-phenylether	10	10 U	10 U	10 U	10 U	10 U
Hexachlorobenzene	10	10 U	10 U	10 U	10 U	10 U
Pentachlorophenol	25	25 U	25 U	25 U	25 U	25 U
Phenanthrene	10	10 U	10 U	10 U	10 U	2 J
Anthracene	10	10 U	10 U	10 U	10 U	10 U
Carbazole	10	10 U	10 U	10 U	10 U	2 J
Di-n-butylphthalate	10	10 U	10 U	10 U	10 U	10 U
Fluoranthene	10	10 U	10 U	10 U	10 U	10 U
Pyrene	10	10 U	10 U	10 U	10 U	10 U
Butylbenzylphthalate	10	10 U	10 U	10 U	10 U	10 U
3,3'-Dichlorobenzidine	10	10 U	10 U	10 U	10 U	10 U
Benzo(a)Anthracene	10	10 U	10 U	10 U	10 U	10 U
Chrysene	10	10 U	10 U	10 U	10 U	10 U
bis(2-Ethylhexyl)phthalate	10	10 U	10 U	2 J	10 U	10 U
Di-n-octylphthalate	10	10 U	10 U	10 U	10 U	10 U
Benzo(b)Fluoranthene	10	10 U	10 U	10 U	10 U	10 U
Benzo(k)Fluoranthene	10	10 U	10 U	10 U	10 U	10 U
Benzo(a)Pyrene	10	10 U	10 U	10 U	10 U	10 U
Indeno(1,2,3-c,d)Pyrene	10	10 U	10 U	10 U	10 U	10 U
Dibenz(a,h)Anthracene	10	10 U	10 U	10 U	10 U	10 U
Benzo(g,h,i)perylene	10	10 U	10 U	10 U	10 U	10 U

Dilution Factor:	1.00	1.00	1.00	1.00
Sample Volume\Weight (ml\g):	1000	1000	1000	1000
Associated Method Blank:	S2097.D	S2097.D	S2097.D	S2097.D
Associated Equipment Blank:	ENQS004XXX94XX	ENQS004XXX94XX	ENQS004XXX94XX	ENQS004XXX94XX
Associated Field Blank:	-	-	-	-

Site: MONITORING WELLS
U: not detected J: estimated

Table 2
Validation / Summary Table

		LOCATION:	CL-102 DUP	CL-102	CL-103	CL-104
		ISIS ID:	ENCL102XXX94XD	ENCL102XXX94XX	ENCL103XXX94XX	ENCL104XXX94XX
		LAB NUMBER:	2223317	2223314	2223318	2223312 D
		DATE SAMPLED:	10/04/94	10/04/94	10/04/94	10/05/94
		DATE EXTRACTED:	10/08/94	10/08/94	10/08/94	10/08/94
		DATE ANALYZED:	10/29/94	10/29/94	10/29/94	11/04/94
ANALYTE	SOW-3/90 - II	CRQL				
alpha-BHC	0.05	0.056 U	0.05 U	0.05 U	0.05 U	0.5 UJ
beta-BHC	0.05	0.056 U	0.05 U	0.05 U	0.05 U	0.5 U
delta-BHC	0.05	0.056 U	0.05 U	0.05 U	0.05 U	0.5 U
gamma-BHC (Lindane)	0.05	0.056 UJ	0.05 UJ	0.05 U	0.05 U	0.5 UJ
Heptachlor	0.05	0.056 UJ	0.05 UJ	0.05 U	0.05 U	0.5 U
Aldrin	0.05	0.056 UJ	0.05 UJ	0.05 U	0.05 U	0.5 U
Heptachlor Epoxide	0.05	0.056 U	0.05 U	0.05 U	0.05 U	0.5 U
Endosulfan I	0.05	0.056 U	0.05 U	0.05 U	0.05 U	0.5 U
Dieldrin	0.1	0.11 UJ	0.1 UJ	0.1 U	0.1 U	1.0 U
4,4'-DDE	0.1	0.11 U	0.1 U	0.1 U	0.1 U	0.99 J
Endrin	0.1	0.11 UJ	0.1 UJ	0.1 U	0.1 U	1.0 U
Endosulfan II	0.1	0.11 U	0.1 U	0.1 U	0.1 U	1.0 U
4,4'-DDD	0.1	0.11 U	0.1 U	0.1 U	0.1 U	1.0 U
Endrin Aldehyde	0.1	0.11 UJ	R	0.1 UJ	0.1 U	1.0 U
Endosulfan Sulfate	0.1	0.11 U	0.1 U	0.1 U	0.1 U	1.0 U
4,4'-DDT	0.1	0.11 UJ	0.1 UJ	0.1 U	0.1 U	1.0 U
Methoxychlor	0.5	0.56 U	0.5 U	0.5 U	0.5 U	5.0 U
Endrin Ketone	0.1	0.11 U	0.1 U	0.1 U	0.1 U	1.0 U
alpha-Chlordane	0.05	0.056 U	0.05 U	0.05 U	0.05 U	0.5 U
gamma-Chlordane	0.05	0.056 U	0.05 U	0.05 U	0.05 U	0.5 U
Toxaphene	5	5.6 U	5.0 U	5.0 U	5.0 U	50 U
Aroclor-1016	1	1.1 U	1.0 U	1.0 U	1.0 U	10 U
Aroclor-1221	2	2.2 U	2.0 U	2.0 U	2.0 U	20 U
Aroclor-1232	1	1.1 U	1.0 U	1.0 U	1.0 U	10 U
Aroclor-1242	1	1.1 U	1.0 U	1.0 U	1.0 U	10 U
Aroclor-1248	1	1.1 U	1.0 U	1.0 U	1.0 U	10 U
Aroclor-1254	1	1.1 U	1.0 U	1.0 U	1.0 U	10 U
Aroclor-1260	1	1.1 U	1.0 U	1.0 U	1.0 U	10 U
=====						
Dilution Factor:		1.00	1.00	1.00	10.0	
Sample Volume/Weight (ml\g):		900	1000	1000	1000	
Associated Method Blank:		PWB1008B	PWB1008B	PWB1008B	PBW1008A2	
Associated Equipment Blank:		ENQSXX2XXX94XX	ENQSXX2XXX94XX	ENQSXX2XXX94XX	ENQSXX2XXX94XX	
Associated Field Blank:		-	-	-	-	

Site: SUMP LIQUIDS

U: not detected J: estimated R: unusable

Table 2
Validation / Summary Table

LOCATION:	CD-101	CD-103
ISIS ID:	ENCD101XXX94XX	ENCD103XXX94XX
LAB NUMBER:	2223303	2223304
DATE SAMPLED:	10/04/94	10/04/94
DATE EXTRACTED:	10/08/94	10/08/94
DATE ANALYZED:	11/13/94	11/13/94

ANALYTE	SOW-3/90 - II	CRQL			
alpha-BHC	1.7		45	JN	R
beta-BHC	1.7			R	R
delta-BHC	1.7			R	R
gamma-BHC (Lindane)	1.7			6.2	J
Heptachlor	1.7			R	R
Aldrin	1.7			R	R
Heptachlor Epoxide	1.7			8.7	J
Endosulfan I	1.7			R	R
Dieldrin	3.3			R	R
4,4'-DDE	3.3			R	R
Endrin	3.3			R	R
Endosulfan II	3.3			R	R
4,4'-DDD	3.3			R	R
Endrin Aldehyde	3.3			R	R
Endosulfan Sulfate	3.3			R	R
4,4'-DDT	3.3			R	R
Methoxychlor	17			R	R
Endrin Ketone	3.3			R	R
alpha-Chlordane	1.7			R	R
gamma-Chlordane	1.7			R	R
Toxaphene	170			R	R
Aroclor-1016	33			R	R
Aroclor-1221	67			R	R
Aroclor-1232	33			R	R
Aroclor-1242	33			R	R
Aroclor-1248	33			R	R
Aroclor-1254	33			R	R
Aroclor-1260	33			R	R

Dilution Factor:	5.00	2.00
Percent Solids:	32	55
Sample Volume\Weight (ml\g):	30.0	30.0
Associated Method Blank:	PSB1008B	PSB1008B
Associated Equipment Blank:	ENQSXX1XXX94XX	ENQSXX1XXX94XX
Associated Field Blank:	-	-

Site: SUMP SEDIMENTS

J: estimated N: presumptive evidence R: unusable

Table 2
Validation / Summary Table

	LOCATION:	WT-101 DUP	WT-101	WT-102	WT-103	WT-104	WT-105	WT-106
	ISIS ID:	ENWT101XXX94XD	ENWT101XXX94XX	ENWT102XXX94XX	ENWT103XXX94XX	ENWT104XXX94XX	ENWT105XXX94XX	ENWT106XXX94XX
	LAB NUMBER:	2223309	2223306	2223305	2223310	2223311	2223301	2223302 D
	DATE SAMPLED:	10/04/94	10/04/94	10/04/94	10/04/94	10/04/94	10/05/94	10/05/94
	DATE EXTRACTED:	10/25/94	10/08/94	10/08/94	10/08/94	10/08/94	10/08/94	10/08/94
	DATE ANALYZED:	12/02/94	11/13/94	11/13/94	11/13/94	11/13/94	11/13/94	11/17/94
ANALYTE	SOW-3/90 - II	CRQL						
alpha-BHC	1.7	2.0 UJ	R	R	29 U	R	R	37 U
beta-BHC	1.7	2.0 UJ	97 U	R	29 U	R	R	R
delta-BHC	1.7	2.0 UJ	97 U	R	29 U	R	R	37 U
gamma-BHC (Lindane)	1.7	R	R	R	R	R	R	37 U
Heptachlor	1.7	R	R	R	29 U	R	3.7 JN	37 U
Aldrin	1.7	R	R	R	48 J	R	R	37 U
Heptachlor Epoxide	1.7	2.0 UJ	R	R	29 U	R	6.2 JN	37 U
Endosulfan I	1.7	R	R	R	96	R	5.9 JN	37 U
Dieldrin	3.3	R	R	R	R	27 JN	3.4 J	73 U
4,4'-DDE	3.3	3.9 UJ	R	50 J	R	51 JN	10 J	73 U
Endrin	3.3	3.9 UJ	190 UJ	R	56 U	R	R	73 U
Endosulfan II	3.3	3.9 UJ	190 U	R	56 U	R	R	73 U
4,4'-DDD	3.3	3.9 UJ	190 U	R	R	R	R	73 U
Endrin Aldehyde	3.3	R	R	R	56 U	R	R	73 U
Endosulfan Sulfate	3.3	3.9 UJ	190 U	R	56 U	R	R	73 U
4,4'-DDT	3.3	7.1 J	R	R	62 JN	R	R	73 U
Methoxychlor	17	20 UJ	970 U	R	290 U	R	R	370 U
Endrin Ketone	3.3	3.9 UJ	190 U	R	56 U	R	R	73 U
alpha-Chlordane	1.7	2.0 UJ	97 U	R	29 U	36 J	4.9 J	37 U
gamma-Chlordane	1.7	2.0 UJ	R	R	32 JN	21 J	R	37 U
Toxaphene	170	200 UJ	9700 U	R	2900 U	R	R	3700 U
Aroclor-1016	33	39 UJ	1900 U	R	560 U	R	R	730 U
Aroclor-1221	67	80 UJ	3800 U	R	1100 U	R	R	1500 U
Aroclor-1232	33	39 UJ	1900 U	R	560 U	R	R	730 U
Aroclor-1242	33	39 UJ	1900 U	R	560 U	R	R	730 U
Aroclor-1248	33	39 UJ	1900 U	R	560 U	R	R	730 U
Aroclor-1254	33	39 UJ	1900 U	R	560 U	R	R	730 U
Aroclor-1260	33	39 UJ	1900 U	R	560 U	R	R	730 U
=====								
Dilution Factor:	1.00	50.0	5.00	10.0	5.00	1.00	20.0	
Percent Solids:	84	88	48	59	60	91	91	
Sample Volume\Weight (ml\g):	30.0	30.0	30.0	30.0	30.0	30.0	30.0	
Associated Method Blank:	PSB1025A	PSB1008B	PSB1008B	PSB1008B	PSB1008B	PSB1008B	PSB1008A	
Associated Equipment Blank:	ENQSXX1XXX94XX	ENQSXX1XXX94XX	ENQSXX1XXX94XX	ENQSXX1XXX94XX	ENQSXX1XXX94XX	ENQSXX1XXX94XX	ENQSXX1XXX94XX	
Associated Field Blank:	-	-	-	-	-	-	-	

Site: WASTE-SLUDGE

U: not detected J: estimated R: unusable N: presumptive evidence

Table 2
Validation / Summary Table

LOCATION:	SS-101	SS-102 DUP	SS-102
ISIS ID:	ENSS101XXX94XX	ENSS102XXX94XD	ENSS102XXX94XX
LAB NUMBER:	2241101	2241105	2241102
DATE SAMPLED:	10/28/94	10/28/94	10/28/94
DATE EXTRACTED:	11/01/94	11/01/94	11/01/94
DATE ANALYZED:	12/08/94	12/08/94	12/08/94

ANALYTE	SOW-3/90 - II	CRQL				
alpha-BHC	1.7		R	2.3	U	2.3 U
beta-BHC	1.7		R	2.3	U	2.3 U
delta-BHC	1.7		R	2.3	UJ	2.3 UJ
gamma-BHC (Lindane)	1.7		R	2.3	U	2.3 U
Heptachlor	1.7	7.3	J	2.3	UJ	2.3 UJ
Aldrin	1.7		R	2.3	U	2.3 U
Heptachlor Epoxide	1.7		R	2.3	U	2.3 U
Endosulfan I	1.7		R	2.3	U	2.3 U
Dieldrin	3.3		R	4.5	UJ	4.5 UJ
4,4'-DDE	3.3		R	4.5	U	4.5 U
Endrin	3.3		R	4.5	U	4.5 U
Endosulfan II	3.3		R		R	4.5 U
4,4'-DDD	3.3		R	4.5	U	4.5 U
Endrin Aldehyde	3.3		R	4.5	U	4.5 U
Endosulfan Sulfate	3.3		R	4.5	U	4.5 U
4,4'-DDT	3.3	14	NJ		R	4.5 U
Methoxychlor	17	94	UJ	23	U	23 UJ
Endrin Ketone	3.3		R		R	R
alpha-Chlordane	1.7		R	2.3	U	2.3 U
gamma-Chlordane	1.7		R	2.3	U	2.3 U
Toxaphene	170		R	230	U	230 U
Aroclor-1016	33		R	45	U	45 U
Aroclor-1221	67		R	91	U	91 U
Aroclor-1232	33		R	45	U	45 U
Aroclor-1242	33		R	45	U	45 U
Aroclor-1248	33		R	45	U	45 U
Aroclor-1254	33		R	45	U	45 U
Aroclor-1260	33		R	45	U	45 U

Dilution Factor:	5.00	1.00	1.00
Percent Solids:	90	74	74
Sample Volume\Weight (ml\g):	30.0	30.0	30.0
Associated Method Blank:	PSB1101A	PSB1101A	PSB1101A
Associated Equipment Blank:	ENQS003XXX94XX	ENQS003XXX94XX	ENQS003XXX94XX
Associated Field Blank:	-	-	-

Site: SURFACE SOILS

U: not detected R: unusable J: estimated N: presumptive evidence

Table 2
Validation / Summary Table

LOCATION:	BS-101	BS-102
ISIS ID:	ENBS101XX694XX	ENBS102XX694XX
LAB NUMBER:	2241106	2241107
DATE SAMPLED:	10/27/94	10/28/94
DATE EXTRACTED:	11/01/94	11/01/94
DATE ANALYZED:	12/08/94	12/08/94

ANALYTE	SOW-3/90 - II	CRQL		
alpha-BHC	1.7	2.2	U	2.1
beta-BHC	1.7	2.2	U	2.1
delta-BHC	1.7	2.2	UJ	2.1
gamma-BHC (Lindane)	1.7	2.2	U	2.1
Heptachlor	1.7	2.2	U	2.1
Aldrin	1.7	2.2	U	2.1
Heptachlor Epoxide	1.7	2.2	U	2.1
Endosulfan I	1.7	2.2	U	2.1
Dieldrin	3.3	4.3	U	4.1
4,4'-DDE	3.3	4.3	U	4.1
Endrin	3.3	4.3	U	4.1
Endosulfan II	3.3	4.3	U	4.1
4,4'-DDD	3.3	4.3	U	4.1
Endrin Aldehyde	3.3	4.3	U	4.1
Endosulfan Sulfate	3.3	4.3	U	4.1
4,4'-DDT	3.3	4.3	U	4.1
Methoxychlor	17	22	U	21
Endrin Ketone	3.3	4.3	U	4.1
alpha-Chlordane	1.7	2.2	U	2.1
gamma-Chlordane	1.7	2.2	U	2.1
Toxaphene	170	220	U	210
Aroclor-1016	33	43	U	41
Aroclor-1221	67	87	U	83
Aroclor-1232	33	43	U	41
Aroclor-1242	33	43	U	41
Aroclor-1248	33	43	U	41
Aroclor-1254	33	43	U	41
Aroclor-1260	33	43	U	35
J				

Dilution Factor:	1.00	1.00
Percent Solids:	77	81
Sample Volume\Weight (ml\g):	30.0	30.0
Associated Method Blank:	PSB1101A	PSB1101A
Associated Equipment Blank:	ENQS003XXX94XX	ENQS003XXX94XX
Associated Field Blank:	-	-

Site: SOIL BORINGS
U: not detected J: estimated

Table 2
Validation / Summary Table

	LOCATION:	MW-101 DUP	MW-101	MW-102	MW-103
	ISIS ID:	ENMW101XXX94XD	ENMW101XXX94XX	ENMW102XXX94XX	ENMW103XXX94XX
	LAB NUMBER:	2263804	2263801	2263805	2263806
	DATE SAMPLED:	11/30/94	11/30/94	11/30/94	11/30/94
	DATE EXTRACTED:	12/03/94	12/03/94	12/03/94	12/03/94
	DATE ANALYZED:	12/12/94	12/12/94	12/12/94	12/12/94
ANALYTE	SOW-3/90 - II	CRQL			
alpha-BHC	0.05	0.05 UJ	0.05 UJ	0.05 UJ	0.05 UJ
beta-BHC	0.05	0.05 U	0.05 U	0.05 U	0.05 UJ
delta-BHC	0.05	0.05 U	0.05 U	0.05 U	0.05 UJ
gamma-BHC (Lindane)	0.05	0.05 U	0.05 U	0.05 U	0.05 UJ
Heptachlor	0.05	0.05 U	0.05 U	0.05 U	0.05 UJ
Aldrin	0.05	0.05 U	0.05 U	0.05 U	0.05 UJ
Heptachlor Epoxide	0.05	0.05 U	0.05 U	0.05 U	0.05 UJ
Endosulfan I	0.05	0.05 U	0.05 U	0.05 U	0.05 UJ
Dieldrin	0.1	0.1 U	0.1 U	0.1 U	0.1 UJ
4,4'-DDE	0.1	0.1 U	0.1 U	0.1 U	0.1 UJ
Endrin	0.1	0.1 U	0.1 U	0.1 U	0.1 UJ
Endosulfan II	0.1	0.1 U	0.1 U	0.1 U	0.1 UJ
4,4'-DDD	0.1	0.1 U	0.1 U	0.1 U	0.1 UJ
Endrin Aldehyde	0.1	0.1 U	0.1 U	0.1 U	0.1 UJ
Endosulfan Sulfate	0.1	0.1 U	0.1 U	0.1 U	0.1 UJ
4,4'-DDT	0.1	0.1 UJ	0.1 UJ	0.1 UJ	0.1 UJ
Methoxychlor	0.5	0.5 U	0.5 U	0.5 U	0.5 UJ
Endrin Ketone	0.1	0.1 U	0.1 U	0.1 U	0.1 UJ
alpha-Chlordane	0.05	0.05 U	0.05 U	0.05 U	0.05 UJ
gamma-Chlordane	0.05	0.05 U	0.05 U	0.05 U	0.05 UJ
Toxaphene	5	5.0 U	5.0 U	5.0 U	5.0 UJ
Aroclor-1016	1	1.0 U	1.0 U	1.0 U	1.0 UJ
Aroclor-1221	2	2.0 U	2.0 U	2.0 U	2.0 UJ
Aroclor-1232	1	1.0 U	1.0 U	1.0 U	1.0 UJ
Aroclor-1242	1	1.0 U	1.0 U	1.0 U	1.0 UJ
Aroclor-1248	1	1.0 U	1.0 U	1.0 U	1.0 UJ
Aroclor-1254	1	1.0 U	1.0 U	1.0 U	1.0 UJ
Aroclor-1260	1	1.0 U	1.0 U	1.0 U	1.0 UJ
	Dilution Factor:	1.00	1.00	1.00	1.00
	Sample Volume\Weight (ml\g):	1000	1000	1000	1000
	Associated Method Blank:	PWB1203A	PWB1203B	PWB1203A	PWB1203A
	Associated Equipment Blank:	ENQS004XXX94XX	ENQS004XXX94XX	ENQS004XXX94XX	ENQS004XXX94XX
	Associated Field Blank:				

Site: MONITORING WELLS

U: not detected J: estimated

Table 2
Validation / Summary Table

LOCATION:	CL-102 DUP	CL-102	CL-103	CL-104
ISIS ID:	ENCL102XXX94XD	ENCL102XXX94XX	ENCL103XXX94XX	ENCL104XXX94XX
LAB NUMBER:	223317	223314	223318	223312
DATE SAMPLED:	10/04/94	10/04/94	10/04/94	10/05/94

ANALYTE	SOW-3/90 - II	CRDL				
Aluminum	200	9040	9520	366	27000	
Antimony	60	38.0 U	38.0 U	38.0 U	38.0 U	
Arsenic	10	5.3 J	5.0 U	5.0 U	54.2	
Barium	200	180 J	203	39.6 J	1320	
Beryllium	5	2.0 U	2.0 U	2.0 U	2.0 U	
Cadmium	5	29.2 J	32.7 J	2.1 J	122 J	
Calcium	5000	53800	58400	68800	580000	
Chromium	10	137	147	9.7 J	1870	
Cobalt	50	30.7 J	38.0 J	21.4 J	375	
Copper	25	1820	1970	120	4930	
Iron	100	70000	91200	20000	265000	
Lead	3	1960	R	R	5510	
Magnesium	5000	10600	12100	18300	57000	
Manganese	15	1300 J	1610 J	2710 J	2330 J	
Mercury	0.2	1.6	2.0	0.20 U*	545	
Nickel	40	155 J	196 J	48.7	3500	
Potassium	5000	4600 J	5350	13200	30800	
Selenium	5	R	R	R	R	
Silver	10	5.0 U	5.0 U	5.0 U	23.1	
Sodium	5000	5840	5870	38300	2030000	
Thallium	10	5.0 U	5.0 U	5.0 UJ	5.0 UJ	
Vanadium	50	67.3	76.5	17.0 U	124	
Zinc	20	7190	8540	337	24500	
Cyanide	10	10.0 U	10.0 U	10.0 U	10.0 U	

Associated Method Blank:	PBENRX1	PBENRX1	PBENRX1	PBENRX1
Associated Equipment Blank:	ENQSXX2XXX94XX	ENQSXX2XXX94XX	ENQSXX2XXX94XX	ENQSXX2XXX94XX
Associated Field Blank:	-	-	-	-

Site: SUMP LIQUIDS

U: not detected J: estimated R: unusable *: duplicate analysis not met

Table 2
Validation / Summary Table

LOCATION:	CD-101	CD-103
ISIS ID:	ENCD101XXX94XX	ENCD103XXX94XX
LAB NUMBER:	223303	223304
DATE SAMPLED:	10/04/94	10/04/94

ANALYTE	SOW-3/90 - II	CRDL		
Aluminum	40	1750	J	4270
Antimony	12	23.1	UJ	16.1 J
Arsenic	2	6.6	J	6.8 J
Barium	40	142	J	186
Beryllium	1	1.2	UJ	0.67 U
Cadmium	1	11.6	J	10.2 J
Calcium	1000	4460	J	17800
Chromium	2	78.4	J	254
Cobalt	10	22.7	J	45.1
Copper	5	181	J	650
Iron	20	66400	J	237000
Lead	0.6	246	J	213
Magnesium	1000	943	UJ	3550
Manganese	3	974	J	1290
Mercury	0.1	1.3	J	1.0
Nickel	8	167	J	406
Potassium	1000	512	UJ	329 J
Selenium	1	3.0	UJ	1.7 UJ
Silver	2	3.0	UJ	1.7 UJ
Sodium	1000	2050	J	360 J
Thallium	2	3.0	UJ	1.7 UJ
Vanadium	10	16.4	J	59.3 J
Zinc	4	1140	J	1140
Cyanide	1	1.5	UJ	0.94 U
=====				
Percent Solids:		32		55

Associated Method Blank:	PBENRX1	PBENRX1
Associated Equipment Blank:	ENQSXX1XXX94XX	ENQSXX1XXX94XX
Associated Field Blank:	-	-

Site: SUMP SEDIMENTS
U: not detected J: estimated

Table 2
Validation / Summary Table

		LOCATION: WT-101 DUP	WT-101	WT-102	WT-103	WT-104	WT-105	WT-106
		ISIS ID: ENWT101XXX94XD	ENWT101XXX94XX	ENWT102XXX94XX	ENWT103XXX94XX	ENWT104XXX94XX	ENWT105XXX94XX	ENWT106XXX94XX
		LAB NUMBER: 223309	223306	223305	223310	223311	223301	223302
		DATE SAMPLED: 10/04/94	10/04/94	10/04/94	10/04/94	10/04/94	10/05/94	10/05/94
ANALYTE	SOW-3/90 - II	CRDL						
Aluminum	40	6370	3910	6380 J	1870	1910	1600	1230
Antimony	12	8.7 UJ	8.3 UJ	13.6 J	12.8 UJ	11.0 UJ	8.3 UJ	16.7 J
Arsenic	2	11.3 J	9.6 J	35.6 J	21.7 J	14.3 J	11.9 J	14.1 J
Barium	40	419	345	184 J	124	110	27.2 J	51.1
Beryllium	1	0.83 J	0.69 J	0.69 UJ	0.67 U	0.58 U	0.44 U	0.38 U
Cadmium	1	36.8 J	21.4 J	9.1 J	18.6 J	8.8 J	4.9 J	10.3 J
Calcium	1000	7110	6720	68000 J	21300	20900	29300	13900
Chromium	2	3730	2790	346 J	102	73.5	26.0	86.9
Cobalt	10	861	912	29.7 J	34.8	26.5	8.6 J	33.3
Copper	5	2770	1800	379 J	553	219	265	296
Iron	20	56900	133000	112000 J	608000	213000	71400	282000
Lead	0.6	1810	1850	363 J	275	46.3	60.5	296
Magnesium	1000	1650	1330	5530 J	2900	2140	4740	2380
Manganese	3	501	916	891 J	2830	2370	407	1360
Mercury	0.1	1.1	1.1	52.9 J	1.8	1.4	0.88	2.3
Nickel	8	11400	9530	209 J	129	78.2	28.3	101
Potassium	1000	1490	1240	1160 J	2460	1460	2170	3400
Selenium	1	5.5 UJ	2.1 J	1.9 UJ	1.5 UJ	1.6 UJ	1.1 UJ	1.0 UJ
Silver	2	16.6 J	20.5 J	1.7 UJ	1.7 UJ	1.5 UJ	1.1 UJ	0.95 UJ
Sodium	1000	6060 J	3680 J	984 J	3050	721 J	6470	8100
Thallium	2	1.1 UJ	1.0 UJ	1.9 UJ	1.5 UJ	1.6 UJ	1.1 UJ	1.0 UJ
Vanadium	10	30.2 J	84.3 J	45.3 J	58.4 J	51.2 J	20.8 J	60.8 J
Zinc	4	2300	2090	1910 J	1370	1530	566	1390
Cyanide	1	0.52 UJ	2.7 J	0.99 UJ	20.7	0.81 U	0.53 U	0.49 U
Percent Solids:			84	88	48	59	60	91
Associated Method Blank:			PBENRX1	PBENRX1	PBENRX1	PBENRX1	PBENRX1	PBENRX1
Associated Equipment Blank:			ENQSXX1XXX94XX	ENQSXX1XXX94XX	ENQSXX1XXX94XX	ENQSXX1XXX94XX	ENQSXX1XXX94XX	ENQSXX1XXX94XX
Associated Field Blank:			-	-	-	-	-	-

Site: WASTE-SLUDGE

U: not detected J: estimated

Table 2
Validation / Summary Table

LOCATION:	SS-101	SS-102 DUP	SS-102
ISIS ID:	ENSS101XXX94XX	ENSS102XXX94XD	ENSS102XXX94XX
LAB NUMBER:	241101	241105	241102
DATE SAMPLED:	10/28/94	10/28/94	10/28/94

ANALYTE	SOW-3/90 - II	CRDL				
Aluminum	40	1650		9860	10200	
Antimony	12	8.0 UJ		9.2 UJ	10.1 UJ	
Arsenic	2	4.4 J		35.1 J	40.5 J	
Barium	40	27.3 J		142	178	
Beryllium	1	0.42 U		4.7	4.0	
Cadmium	1	1.8 J		1.1 J	2.3 J	
Calcium	1000	212000		34900	35900	
Chromium	2	11.4		10.1	14.3	
Cobalt	10	6.1 J		10.9 J	12.1 J	
Copper	5	67.9		77.9	69.4	
Iron	20	7870		12800	18200	
Lead	0.6	R		R	R	
Magnesium	1000	18300		7060	6400	
Manganese	3	495		380	284	
Mercury	0.1	0.40		0.61	0.88	
Nickel	8	24.5		25.9	30.8	
Potassium	1000	470 J		521 J	1010 J	
Selenium	1	1.1 UJ		1.3 U	1.3 U	
Silver	2	1.0 U		1.2 U	1.3 U	
Sodium	1000	106 J		237 J	190 J	
Thallium	2	1.1 UJ		1.3 U	1.3 U	
Vanadium	10	11.7		34.4	35.4	
Zinc	4	653		2100	1940	
Cyanide	1	0.53 U		0.57 U	0.58 U	
=====						
Percent Solids:		90		74	74	

Associated Method Blank:	PBENRX2	PBENRX2	PBENRX2
Associated Equipment Blank:	ENQS003XXX94XX	ENQS003XXX94XX	ENQS003XXX94XX
Associated Field Blank:	-	-	-

Site: SURFACE SOILS

U: not detected J: estimated R: unusable

Table 2
Validation / Summary Table

LOCATION:	BS-101	BS-102
ISIS ID:	ENBS101XX694XX	ENBS102XX694XX
LAB NUMBER:	241106	241107
DATE SAMPLED:	10/27/94	10/28/94

ANALYTE	SOW-3/90 - 11	CRDL		
Aluminum	40	7130		16100
Antimony	12	9.2 UJ		9.4 J
Arsenic	2	5.1 J		4.6 J
Barium	40	57.0		186
Beryllium	1	0.48 U		0.93 J
Cadmium	1	0.48 UJ		0.47 UJ
Calcium	1000	12200		19700
Chromium	2	8.7		20.8
Cobalt	10	6.4 J		10.2 J
Copper	5	49.5		25.9
Iron	20	15600		31100
Lead	0.6	R		R
Magnesium	1000	4780		8670
Manganese	3	270		485
Mercury	0.1	0.13 U		0.12 U
Nickel	8	12.9		26.3
Potassium	1000	982 J		2470
Selenium	1	1.1 UJ		1.2 U
Silver	2	1.2 U		1.2 U
Sodium	1000	1330		295 J
Thallium	2	1.1 UJ		1.2 U
Vanadium	10	19.9		42.6
Zinc	4	65.4		117
Cyanide	1	0.67 U		0.52 U
=====				
Percent Solids:		77		81

Associated Method Blank:	PBENRX2	PBENRX2
Associated Equipment Blank:	ENQS003XXX94XX	ENQS003XXX94XX
Associated Field Blank:	-	-

Site: SOIL BORINGS

U: not detected J: estimated R: unusable

Table 2
Validation / Summary Table

LOCATION:	MW-101 DUP	MW-101	MW-102	MW-103
ISIS ID:	ENMW101XXX94XD	ENMW101XXX94XX	ENMW102XXX94XX	ENMW103XXX94XX
LAB NUMBER:	263804	263801	263805	263806
DATE SAMPLED:	11/30/94	11/30/94	11/30/94	11/30/94

ANALYTE	SOW-3/90 - 11	CRDL				
Aluminum	200	2390	J	250	J	315
Antimony	60	52.9	J	38.0	UJ	38.0
Arsenic	10	5.0	U	5.0	U	5.0
Barium	200	88.7	J	82.5	J	80.4
Beryllium	5	2.0	U	2.0	U	2.0
Cadmium	5	2.0	UJ	2.0	UJ	2.0
Calcium	5000	136000		128000		212000
Chromium	10	7.2	J	5.0	UJ	5.0
Cobalt	50	6.0	U	6.0	U	6.0
Copper	25	11.8	J	5.0	U	5.4
Iron	100	7760		6240		1850
Lead	3	11.2	J	3.0	UJ	15.7
Magnesium	5000	54600		53100		41000
Manganese	15	412		339		2990
Mercury	0.2	0.20	U	0.20	U	0.20
Nickel	40	26.0	U	26.0	U	26.0
Potassium	5000	14200	J	14200	J	49200
Selenium	5	5.0	U	5.0	U	5.0
Silver	10	5.0	U	5.0	U	5.0
Sodium	5000	107000		110000		208000
Thallium	10	5.0	U	5.0	U	5.0
Vanadium	50	17.0	U	17.0	U	17.0
Zinc	20	13.6	J	5.0	U	24.7
Cyanide	10	10.0	U	10.0	U	10.0

Associated Method Blank:	PBENRX3	PBENRX3	PBENRX3	PBENRX3
Associated Equipment Blank:	ENQS004XXX94XX	ENQS004XXX94XX	ENQS004XXX94XX	ENQS004XXX94XX
Associated Field Blank:	-	-	-	-

Site: MONITORING WELLS
U: not detected J: estimated

Table 2
Validation / Summary Table

LOCATION:		CL-103	CL-104
ISIS ID:		ENCL103XXX94XX	ENCL104XXX94XX
LAB NUMBER:		D223318	D223312
DATE SAMPLED:		10/04/94	10/05/94
ANALYTE	RL		
arsenic	52.0/5.0	5.0 U	10.0 UJ
barium	11.0	R	R
cadmium	2.0	3.8 J	2.0 U
chromium	5.0	7.9 J	5.0 UJ
lead	26.0/3.0	94.3 J	146 J
mercury	0.20	0.20 U	0.57
selenium	90.0/5.0	5.0 UJ	10.0 UJ
silver	5.0	5.0 UJ	5.0 UJ

Associated Method Blank: EPENRX1 EPENRX1
Associated Equipment Blank: ENQSXX2XXX94XX ENQSXX2XXX94XX
Associated Field Blank: - -

SITE: SUMP LIQUIDS

Note: Inorganic Data - EPTOX Metals

U: not detected R: unusable J: estimated

Table 2
Validation / Summary Table

	LOCATION:	CD-101	CD-103
	ISIS ID:	ENCD101XXX94XX	ENCD103XXX94XX
	LAB NUMBER:	E223303	E223304
	DATE SAMPLED:	10/04/94	10/04/94

ANALYTE	RL		
arsenic	52.0/5.0	52.0 U	52.0 U
barium	11.0	R	R
cadmium	2.0	2.0 U	2.0 U
chromium	5.0	5.0 UJ	5.0 UJ
lead	26.0/3.0	26.0 UJ	26.0 UJ
mercury	0.20	0.20 U	0.20 U
selenium	90.0/5.0	90.0 U	90.0 U
silver	5.0	5.0 UJ	5.0 UJ

Associated Method Blank: EPENRX1 EPENRX1
Associated Equipment Blank: ENQSXX1XXX94XX ENQSXX1XXX94XX
Associated Field Blank: - -

SITE: SUMP SEDIMENTS

Note: Inorganic Data - EPTOX Metals

U: not detected R: unusable J: estimated

Table 2
Validation / Summary Table

ANALYTE	WT-101 DUP		WT-101		WT-102		WT-103		WT-104		WT-105		WT-106	
	LOCATION: WT-101 DUP		WT-101		WT-102		WT-103		WT-104		WT-105		WT-106	
	ISIS ID: ENWT101XXX94XD		ENWT101XXX94XX		ENWT102XXX94XX		ENWT103XXX94XX		ENWT104XXX94XX		ENWT105XXX94XX		ENWT106XXX94XX	
	LAB NUMBER: E223309		E223306		E223305		E223310		E223311		E223301		E223302	
DATE SAMPLED:		10/04/94	10/04/94		10/04/94		10/04/94		10/04/94		10/05/94		10/05/94	
RL														
arsenic	52.0/5.0	52.0 U	52.0 U		52.0 U		52.0 U		52.0 U		52.0 U		52.0 U	
barium	11.0	R	R		R		R		R		R		R	
cadmium	2.0	1100	1100		3.7 J		22.5		2.0 U		17.0 J		10.9 J	
chromium	5.0	84.3 J	51.0 J		5.9 J		9.5 J		5.0 UJ		53.6 J		25.0 J	
lead	26.0/3.0	4820 J	11400 J		26.0 UJ		58.1 J		26.0 UJ		45.2 J		60.8 J	
mercury	0.20	0.25	0.34		0.20 U		0.23		0.20 U		0.52		1.2	
selenium	90.0/5.0	90.0 U	90.0 U		90.0 U		90.0 U		90.0 U		90.0 U		90.0 U	
silver	5.0	37.6 J	33.4 J		5.0 UJ		5.0 UJ		5.0 UJ		5.0 UJ		5.0 UJ	

=====

Associated Method Blank:	EPENRX1	EPENRX1	EPENRX1	EPENRX1	EPENRX1	EPENRX1	EPENRX1
Associated Equipment Blank:	ENQSXX1XXX94XX	ENQSXX1XXX94XX	ENQSXX1XXX94XX	ENQSXX1XXX94XX	ENQSXX1XXX94XX	ENQSXX1XXX94XX	ENQSXX1XXX94XX
Associated Field Blank:	-	-	-	-	-	-	-

SITE: WASTE-SLUDGE

Note: Inorganic Data - EPTOX Metals

U: not detected R: unusable J: estimated

Table 2
Validation / Summary Table

	LOCATION:	SS-101	SS-102 DUP	SS-102
	ISIS ID:	ENSS101XXX94XX	ENSS102XXX94XD	ENSS102XXX94XX
	LAB NUMBER:	E241101	E241105	E241102
	DATE SAMPLED:	10/28/94	10/28/94	10/28/94
ANALYTE	RL			
arsenic	52.0	52.0 U	52.0 U	52.0 U
barium	11.0	507 J	1000 J	1080 J
cadmium	2.0	2.0 UJ	15.8 J	15.8 J
chromium	5.0	5.0 UJ	6.0 J	5.0 UJ
lead	26.0	26.0 U	33.7 J	33.7 J
mercury	0.2	0.2 U	0.2 U	0.2 U
selenium	90.0	90.0 U	90.0 U	90.0 U
silver	5.0	5.0 U	5.0 U	5.0 U

Associated Method Blank: EPENRX2 EPENRX2 EPENRX2
Associated Equipment Blank: - - -
Associated Field Blank: - - -

SITE: SURFACE SOILS
Note: Inorganic Data - EPTOX Metals
U: not detected J: estimated

Table 2
Validation / Summary Table

LOCATION:	CL-103	CL-104
ISIS ID:	ENCL103XXX94XX	ENCL104XXX94XX
LAB NUMBER:	2223318	2223312
DATE SAMPLED:	10/04/94	10/05/94
DATE ANALYZED:	10/12/94	10/12/94

ANALYTE	RL		
Corrosivity, inch/Year	0.01	0.01 U	0.01 U
Ignitability, Degrees F		>212	>212
Cyanide, Reactive, ppm	1	1 U	1 U
Sulfide, Reactive, ppm	1	1 U	1 U

=====

Associated Method Blank:	WCENRX1	WCENRX1
Associated Equipment Blank:	ENQSXX2XXX94XX	ENQSXX2XXX94XX
Associated Field Blank:	-	-

SITE: SUMP LIQUIDS
U: not detected

Table 2
Validation / Summary Table

LOCATION:	CD-101	CD-103
ISIS ID:	ENCD101XXX94XX	ENCD103XXX94XX
LAB NUMBER:	2223303	2223304
DATE SAMPLED:	10/04/94	10/04/94
DATE ANALYZED:	10/12/94	10/12/94

ANALYTE	RL		
Corrosivity, inch/Year	0.01	0.01 U	0.01 U
Ignitability, Degrees F		>212	>212
Cyanide, Reactive, ppm	1	1 U	1 U
Sulfide, Reactive, ppm	1	1 U	1 U

=====

Associated Method Blank:	WCENRX1	WCENRX1
Associated Equipment Blank:	ENQSXX1XXX94XX	ENQSXX1XXX94XX
Associated Field Blank:	-	-

SITE: SUMP SEDIMENTS
U: not detected

Table 2
Validation / Summary Table

ANALYTE	RL	LOCATION: WT-101 DUP	WT-101	WT-102	WT-103	WT-104	WT-105	WT-106
		ISIS ID: ENWT101XXX94XD	ENWT101XXX94XX	ENWT102XXX94XX	ENWT103XXX94XX	ENWT104XXX94XX	ENWT105XXX94XX	ENWT106XXX94XX
		LAB NUMBER: 2223309	2223306	2223305	2223310	2223311	2223301	2223302
		DATE SAMPLED: 10/04/94	10/04/94	10/04/94	10/04/94	10/04/94	10/05/94	10/05/94
		DATE ANALYZED: 10/12/94	10/12/94	10/12/94	10/12/94	10/12/94	10/12/94	10/12/94
Corrosivity, inch/Year	0.01	0.01 U	0.01 U	0.01 U	0.01 U	0.01 U	0.01 U	0.01 U
Ignitability, Degrees F		>212	>212	>212	>212	>212	>212	>212
Cyanide, Reactive, ppm	1	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Sulfide, Reactive, ppm	1	1 U	1 U	1 U	1 U	1 U	1 U	1 U

```

=====
Associated Method Blank:  WCENRX1      WCENRX1      WCENRX1      WCENRX1      WCENRX1      WCENRX1      WCENRX1
Associated Equipment Blank: ENQSXX1XXX94XX ENQSXX1XXX94XX ENQSXX1XXX94XX ENQSXX1XXX94XX ENQSXX1XXX94XX ENQSXX1XXX94XX ENQSXX1XXX94XX
Associated Field Blank:      -            -            -            -            -            -            -

```

SITE: WASTE-SLUDGE

U: not detected

Table 2
Validation / Summary Table

	LOCATION:	SS-101	SS-102 DUP	SS-102
	ISIS ID:	ENSS101XXX94XX	ENSS102XXX94XD	ENSS102XXX94XX
	LAB NUMBER:	2241101	2241105	2241102
	DATE SAMPLED:	10/28/94	10/28/94	10/28/94
	DATE ANALYZED:	11/14/94	11/14/94	11/14/94

ANALYTE	RL			
Corrosivity, inch/Year	0.01	0.01 U	0.01 U	0.01 U
Ignitability, Degrees f		>212	>212	>212
Cyanide, Reactive, ppm	1	1 U	1 U	1 U
Sulfide, Reactive, ppm	1	1 U	1 U	1 U

=====

Associated Method Blank:	WCENRX2	WCENRX2	WCENRX2
Associated Equipment Blank:	ENQS003XXX94XX	ENQS003XXX94XX	ENQS003XXX94XX
Associated Field Blank:			

SITE: SURFACE SOILS
U: not detected

TENTATIVELY IDENTIFIED COMPOUND (TIC) SUMMARY
 NYSDEC-PSA-14 ENRX Inc. Site; FILE: 7170-01
 SOIL (ug/kg)

VOLATILE

	ENWT101XXX94XD	ENWT101XXX94XX	ENWT101XXX94XXDL	ENCD101XXX94XX
unknown aromatic	5600 J(3)	11000 J(3)	14000 J	2000 J
unknown	800 J			
trimethyl benzene isomer	5000 J(2)	10000 J(3)	7900 J	
ethyl dimethyl benzene isomer	780 J	2300 J(2)		
methyl propyl benzene isomer	780 J	1300 J		
ethyl methyl benzene isomer	1100 J	2500 J	4400 J	
unknown hydrocarbon				20000 J(9)

	ENWT102XXX94XX	ENWT106XXX94XX
unknown hydrocarbon	63 J(4)	
unknown		8 J

NO VOLATILE TIC'S WERE IDENTIFIED IN THE FOLLOWING SAMPLES:

ENCD103XXX94XX
 ENWT103XXX94XX
 ENWT104XXX94XX
 ENWT105XXX94XX

SEMIVOLATILE

	ENCD101XXX94XX	ENCD103XXX94XX	ENWT101XXX94XX	ENWT101XXX94XD
unknown hydrocarbon	7500 J		2000000 J(17)	63000 J(4)
unknown	360000 J(4)	86000 J(6)	250000 J(2)	170000 J(11)
unknown aromatic				54000 J(4)

	ENWT102XXX94XX	ENWT103XXX94XX	ENWT104XXX94XX	ENWT105XXX94XX
unknown hydrocarbon	56000 J(3)	33000 J(2)	160000 J(18)	19000 J(3)
unknown	490000 J(12)	140000 J(17)	18000 J(2)	73000 J(6)
unknown aromatic	38000 J(2)	7000 J(2)		20000 J(2)

	ENWT106XXX94XX
unknown	190000 J(12)

Data Qualifiers: J - estimated

TENTATIVELY IDENTIFIED COMPOUND (TIC) SUMMARY
 NYSDEC-PSA-14 ENRX Inc. Site; FILE: 7170-01
 AQUEOUS (ug/L)

VOLATILE

	ENCL102XXX94XD	ENCL102XXX94XX	ENCL104XXX94XX	ENCL104XXX94XXDL
unknown aromatic	5 J		370 J	
unknown freon isomer		6 J	79000 J	83000 J
naphthalene isomer		16 J		
unknown			1300 J(3)	
unknown hydrocarbon			180 J	

NO VOLATILE TIC'S WERE IDENTIFIED IN THE FOLLOWING SAMPLES:

ENCL102XXX94XDDL ENCL102XXX94XXDL
 ENCL103XXX94XX

SEMIVOLATILE

	ENQSXX1XXX94XX	ENCL102XXX94XX	ENCL102XXX94XD	ENCL103XXX94XX
unknown	120 J(10)	2000 J(4)	1200 J(5)	300 J(5)
unknown siloxane	8 J(2)			
unknown hydrocarbon		300 J(4)	640 J(11)	69 J(4)
unknown aromatic			47 J	21 J

ENCL104XXX94XXDL

unknown	5700 J(8)
unknown hydrocarbon	4700 J(7)
unknown aromatic	820 J(2)

Data Qualifiers: J - estimated

TENTATIVELY IDENTIFIED COMPOUND (TIC) SUMMARY
NYSDEC-PSA-14 ENRX Inc. Site; FILE: 7170-02
AQUEOUS (ug/L)

VOLATILE

NO VOLATILE TIC'S WERE IDENTIFIED IN THE FOLLOWING SAMPLES:

ENQS003XXX94XX

TENTATIVELY IDENTIFIED COMPOUND (TIC) SUMMARY
NYSDEC-PSA-14 ENRX Inc. Site; FILE: 7170-02
SOIL (ug/kg)

VOLATILE

NO VOLATILE TIC's WERE IDENTIFIED IN THE FOLLOWING SAMPLES:

ENBS101XX694XX
ENBS102XX694XX
ENSS102XXX94XD
ENSS101XXX94XX
ENSS102XXX94XX

TENTATIVELY IDENTIFIED COMPOUND (TIC) SUMMARY
NYSDEC-PSA-14 ENRX Inc. Site; FILE: 7170-03
AQUEOUS (ug/L)

SEMIVOLATILE

ENQS003XXX94XX	
unknown	J(5)

Data Qualifiers: J - estimated

TENTATIVELY IDENTIFIED COMPOUND (TIC) SUMMARY
 NYSDEC-PSA-14 ENRX Inc. Site; FILE: 7170-03
 SOIL (ug/kg)

SEMIVOLATILE

	ENBS101XX694XX	ENBS102XX694XX	ENSS101XXX94XX	ENSS102XXX94XX
unknown hydrocarbon	450 J(2)	87 J		190 J
unknown aromatic		94 J	12000 J(13)	1000 J(4)
unknown			4300 J(5)	
ENSS102XXX94XD				
unknown aromatic	180 J			

Data Qualifiers: J - estimated

TENTATIVELY IDENTIFIED COMPOUND (TIC) SUMMARY
 NYSDEC-PSA-14 ENRX Inc. Site; FILE: 7170-04
 AQUEOUS (ug/L)

VOLATILE

ENMW103XXX94XX	
unknown	80 J(4)
unknown aromatic	9 J

NO VOLATILE TIC'S WERE IDENTIFIED IN THE FOLLOWING SAMPLES:

ENMW101XXX94XX	ENQS004XXX94XX
ENMW101XXX94XD	ENQT002XXX94XX
ENMW102XXX94XX	

SEMIVOLATILE

	ENMW101XXX94XD	ENMW102XXX94XX	ENMW103XXX94XX	ENQS004XXX94XX
unknown	2 J	49 J(5)	1000 J(7)	100 J(5)
unknown hydrocarbon		89 J(14)		

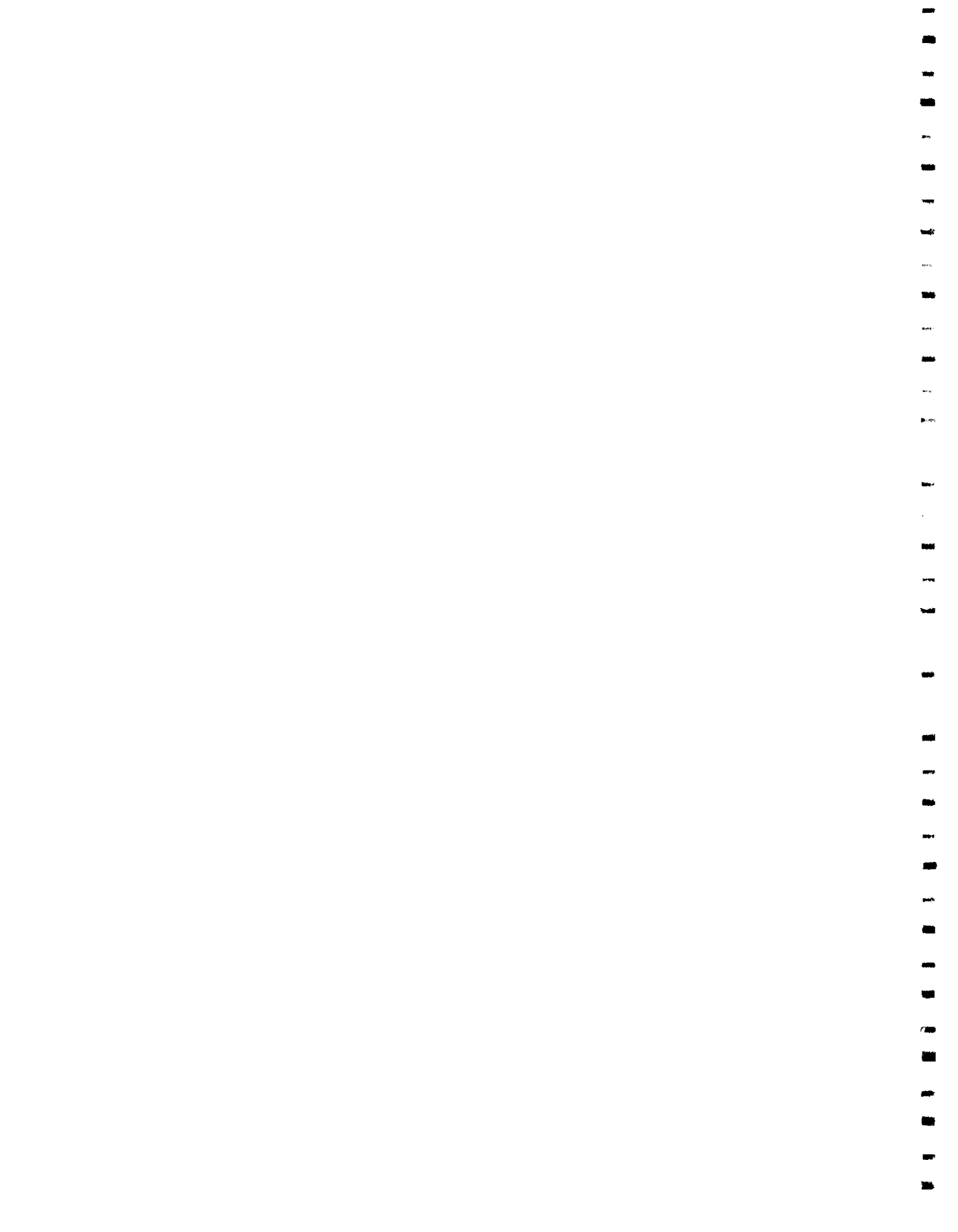
NO SEMIVOLATILE TIC'S WERE IDENTIFIED IN THE FOLLOWING SAMPLES:

ENMW101XXX94XX

Data Qualifiers: J - estimated

SECTION 4.0
SURVEY CONTROL REPORT

ABB Environmental Services



New York State Department of Environmental Conservation

SUPERFUND STANDBY CONTRACT

ENRX INC.

Buffalo, New York

CONTROL REPORT

January 1995



OM P. POPLI, P.E., L.S., P.C.
Consulting Engineers & Surveyors
44 Saginaw Drive
Rochester, NY 14623
(716) 442-6940



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ATTACHMENT A

INTRODUCTION



INTRODUCTION

This report summarizes the results of a site investigation survey encompassing approximately 0.5 acre surrounding the ENRX Inc. located in Buffalo, New York.

The survey work needed to satisfy the requirements of the Task Order Memorandum was performed in December, 1994 by Om P. Popli, P.E., L.S., P.C., by Brad Lins, Party Chief, under the supervision of Michael F. Ives, P.L.S.

ATTACHMENT B
STATEMENT OF WORK

TASK ORDER MEMORANDUM ____

SURVEYING AND MAPPING PRELIMINARY SITE ASSESSMENTS

The services to be provided under Task Order Memorandum ____ shall be performed in accordance with the terms and conditions of the Task Order Agreement between Om Popli Associates Incorporated (POPLI) and ABB Environmental Services (ABB-ES) dated May 5, 1991.

PROJECT SUMMARY

ABB-ES, under contract to the New York State Department of Environmental Conservation (NYSDEC), is performing Preliminary Site Assessments (PSAs) of suspected inactive hazardous waste sites in the State of New York. The purpose of the investigation is to confirm or deny the presence of hazardous waste disposal on-site and determine if a significant threat exists to public health and the environment. Task 1 activities include a data and records search and a site walkover. Task 2 involves the preparation of Work Plans for additional site investigations. Tasks 3 and 4 include initial environmental sampling and subsurface investigations, respectively.

Tasks 1 and 2 been completed and ABB-ES is developing site-specific budgets for the field investigation activities to be conducted under Tasks 3 and 4. As part of Tasks 3 and 4 the services of a licensed land surveyor are required to locate exploration locations, site features, and prepare a map for each site.

SCOPE OF SERVICES

POPLI shall provide all necessary personnel, equipment, and materials to perform the following Scope of Services in accordance with the Standard Specification described in Attachment A.

The follwoing seven sites are included under this Task Order Memorandum are:

SITE LIST

SITE NAME	NYSDEC SITE NO.	CITY/TOWN	COUNTY
Wantagh Cleaners	130054	Hempstead	Nassau
Ranco Wiping Cloth	130076	Freeport	Nassau
Green Thumb Spray Company	130518	Hempstead	Nassau
Target Products, Inc.	819015	LeRoy	Genessee

SITE LIST

SITE NAME	NYSDEC SITE No.	CITY/TOWN	COUNTY
(continued)			
Davidson's Collision	828091	Rochester	Monroe
Hanna Furnace, Division of National Steel Corp	915029	Buffalo	Erie
ENRX, Inc.	915150	Buffalo	Erie

A site location map and site sketch for each site are provided in Attachment B.

GENERAL SERVICES. The following general services are to be provided at each of the seven sites.

1. Mobilize and demobilize all necessary survey equipment and personnel to complete the horizontal location and vertical elevation survey within the project schedule.
2. Supply POPLI's personnel with all necessary equipment and clothing including, but not limited to, hardhats and safety glasses and other items in addition to those normally utilized by POPLI at a nonhazardous site.
3. Maintain good relations with NYSDEC, the local community, and associated agencies and land owners. POPLI field personnel employed on the project should be made thoroughly cognizant of the importance of this aspect of the work and its sensitivity to the entire program.
4. Attend a health and safety meeting with ABB-ES and the NYSDEC prior to the start of the survey activities.
5. Establish appropriate horizontal and vertical control at the site (i.e., locating existing benchmarks) Refer to the Attachment A, Technical Specifications for appropriate control.
6. Establish horizontal control at all monitoring wells, borings, sample locations, corners of buildings, and other points as determined by ABB-ES and indicate on map. Horizontal positions will be tied into the New York State Plane Coordinate System. Horizontal accuracy is to be 0.1 foot.
7. Establish vertical control at all monitoring wells, borings, sample locations, and corners of buildings as determined by ABB-ES and indicate on map. Vertical

elevations will be tied to mean sea level, 1929 General Adjustment. Vertical elevation accuracy will be 0.01 foot.

8. Locate and indicate specific features of the site on the map, such as the location and extent of filled areas, buried tanks, waste piles, buildings, etc. as determined by ABB-ES.
9. Provide all necessary measures for securing POPLI's equipment during the conduct of the work.
10. Conduct all field activities in an efficient and professional manner with minimum impact to the site environment. Tree and brush removal and other activities which impact the existing site environment shall not be undertaken without prior approval by ABB-ES.
11. Prepare a map showing property and site boundaries, developed through the use of current tax maps. The name of current property owners are to be shown on the map. In addition the map shall contain north arrow, scale, a legend that shows designations (wells, borings, sample locations, etc.) and a title block containing the official site name and site number.
12. Provide an electronic copy of the map on a 3.5-inch diskette in a format compatible with AutoCADD Release 12.
13. Provide a final bound report for each site summarizing coordinates of all surveyed locations, and ground elevations, together with any comments pertinent to each location. Sampling locations shall be referenced by identification numbers, to be provided by ABB-ES. This report shall also contain photocopies of all field notes and calculations as an appendix. The report shall describe procedures, traverses, and closures, and will note any significant observations relative to the survey. The final report shall be complete and accurate and shall not contain any errors. Any errors or omissions by POPLI shall be corrected by POPLI at no cost to ABB-ES within two weeks of notice of errors/omissions, so as not to jeopardize the overall project schedule. The final report shall be signed by a surveyor licensed in the State of New York.
14. Provide current health and safety certificates of all POPLI field personnel assigned to field surveying activities at any of the seven sites.

The methods, procedures and techniques to be used by POPLI are the responsibility of POPLI, and shall be designed to meet the intent of the specifications in Attachment A, appended hereto and incorporated by this Task Order Memorandum. Should the technical specifications conflict in any manner with the scope of services, the provisions of the scope of services shall govern.

SITE-SPECIFIC SERVICES. Specific requirements for each site are as follows:

Wantagh Cleaners

Map the site, located at the corners of Wantagh Avenue and Sandhill Road in Hempstead, New York. Include the following items in the survey:

- horizontal locations of 3 monitoring wells;
- vertical elevations of monitoring wells including top of the riser, top of the protective casing, and the ground surface;
- major site characteristics including the building, paved areas, and leaching pools (as indicated by manhole covers);
- property boundaries based on tax map data; and
- 10 miscellaneous spot locations to be established by ABB-ES.

Ranco Wiping Cloth

Map the approximately 0.24-acre site located at 409 North Main Street, Freeport, New York. Include the following items in the survey:

- horizontal locations of 5 shallow subsurface soil samples;
- horizontal locations of 3 monitoring wells;
- vertical elevations of monitoring wells including top of the riser, top of the protective casing, and the ground surface;
- major site characteristics including the edge of paved areas, building corners, and the drywell;
- property boundaries based on tax map data; and
- 10 miscellaneous spot locations to be established by ABB-ES.

Green Thumb Spray Company

Map the approximately 0.2-acre site located at 627 Peninsula Boulevard, Hempstead, New York. Include the following items in the survey:

- horizontal locations of up to 3 shallow subsurface soil samples;
- horizontal locations of 3 monitoring wells;
- vertical elevations of monitoring wells including top of the riser, top of the protective casing, and the ground surface;
- major site characteristics including edge of the paved area, building corners, and drywell;
- property boundaries based on tax map data; and
- 10 miscellaneous spot locations to be established by ABB-ES.

- ~~vertical elevations of monitoring wells including top of the riser, tope of the protective casing, and the ground surface;~~
- ~~major site characteristics including the outline of the Union Ship Canal, existing roads, and existing buildings;~~
- ~~property boundaries based on tax map data; and~~
- ~~50 miscellaneous spot locations to be established by ABB-ES.~~

ENRX, Inc.

Map the approximately 0.5-acre site located at 766 Babcock Street, Buffalo, New York. Include the following items in the survey:

- horizontal locations of 3 monitoring wells;
- vertical elevations of monitoring wells including top of the riser, tope of the protective casing, and the ground surface;
- major site characteristics including edge of paved area, building corners, fenced areas, and locations of utility manholes;
- property boundaries based on tax map data; and
- 10 miscellaneous spot locations to be established by ABB-ES.

ABB-ES or its designated representative will provide the following services:

1. Provide POPLI with right-of-access to all locations through NYSDEC and appropriate land owners.
2. Conduct health and safety meeting with POPLI field representatives prior to initiation of survey activities.

HEALTH AND SAFETY REQUIREMENTS

Before field work begins, POPLI must submit certification to ABB-ES that each of its employees working on-site at the PSA sites is in a Medical Monitoring program and has completed the appropriate training and field experience in compliance with the new OSHA 29 CFR regulations.

POPLI is responsible for meeting the requirements of the laws and regulations that apply to its work and to its employees. POPLI is advised to investigate the new requirements of 29 CFR before beginning work on this project. All work will be done at Level D, as described in the Health and Safety Plan (HASP) which will be forwarded to POPLI by ABB-ES prior to authorization to proceed for individual sites.

PROJECT SCHEDULE

Each site survey will be schedule separately depnding on field work schedules. POPLI shall mobilize within five (5) calendar days of notice to proceed. ABB-ES anticipates that the survey for Tasks 3 and 4 field investigation activities shall commence on or about September 1, 1994. The final bound report shall be completed and provided to ABB-ES no later than 21 calender days after completion of the field work.

NYSDEC or ABB-ES reserve the right to reduce or increase the number of sampling locations, spot elevations, or temporary bench marks in this program. POPLI will provide sufficient equipment and manpower to avoid unnecessary delays.

POPLI shall assume 8-hour days and a repeating schedule of normal 5-day work week with 2 days off on weekends. If survey work falls behind schedule, POPLI shall be prepared to work reasonable overtime and mobilize additional survey equipment and personnel to complete the program within the project schedule, as specified by ABB-ES.

MEASUREMENT AND PAYMENT

The payment items shall be those presented in accordance with contract required rates included in the Contract Schedules included as Attachment C. All measurements for payment purposes shall be rounded to the nearest 0.5 hour and half day. All unit cost shall be based on "Level D Protection". Prevailing Rates do not apply to investigative activities in the PSA. A separate rate schedule is to be provided for each of the fourteen sites.

COMPENSATION

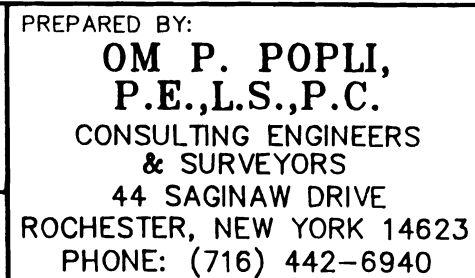
POPLI shall be compensated on a cost plus fixed fee basis for the services described in the Scope of Services as authorized and accepted by ABB-ES in accordance with the contract required rates Attachment C and the site-specific estimates included in Attachment D. Invoices shall include the project job number task order number and indicate by date, the hours, expenses, and services provided on a site by site basis. Time reports, expense reports, and itemization of miscellaneous charges shall be required as backup for each submitted invoice in accordance with the Task Order Agreement-Attachment A: Schedule B, Payment Requirements.

ATTACHMENT C

FIGURES

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1. The first part of the document is a list of names and dates, which appears to be a record of some kind. The names are written in a cursive script, and the dates are in a standard font. The list is organized into two columns, with names on the left and dates on the right. The names are: John Smith, James Brown, William Jones, Thomas White, Robert Black, and Charles Green. The dates are: 1810, 1811, 1812, 1813, 1814, and 1815. The list is followed by a signature and a date.



ENRX INC.

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85-34-21	89-49-38	48.09	-	"
86-11-17	89-44-09	34.60	-	"
08-35-56	TAPED	5.28	-	CLF
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76-44-55	90-06-00	96.04	"	MW-102
76-43-24	89-56-52	113.45	"	BLD
81-04-19	88-05-20	111.81	9.00	EP
80-36-36	87-56-13	98.35	"	"
78-47-51	88-22-13	73.51	7.70	EP
88-41-15	76-48-14	48.92	6.70	"

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ENRX INC.

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	92-22-31	90-40-41	37.87	"	BLD	JOG
	166-17-46	91-18-23	6.48	"	FH	SITE BM HYD
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D	01-05-06	85-27-39	17.17	"	BLD	CNR
N	101-54-26	89-30-35	57.37	"	"	"
	358-49-35	89-18-38	129.70	"	BLD	CNR

ENRX INC.

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BS	D	180-00-00	90-16-29	224.440	30
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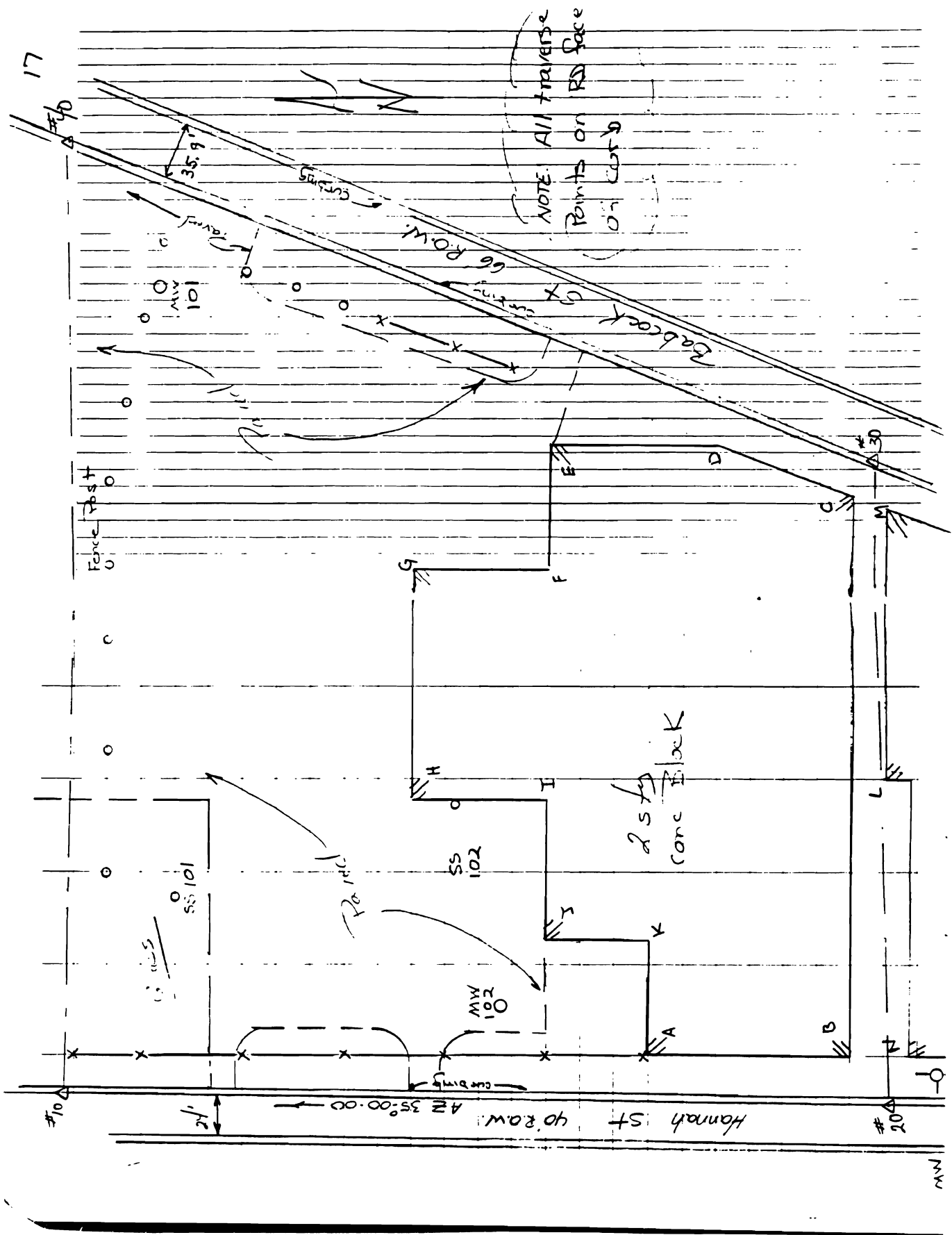
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 FS D 297-13-26 89-48-02 224.430 40
 FS ID 117-13-49 270-11-58 224.430 40
 BS ID 0-00-23 270-35-44 134.335 20



12/22/74
BL TF
Sunny 50°

ENRX INC. HAZ. SITE

CITY OF BUFFALO -

$\pi \approx 20$ BS ≈ 10 FS ≈ 30
HI = 5.24 HR = 5.23 HR = 4.93

BS	D	0-00-00	89-45-22	218.335	10
FS	D	89-53-45	90-43-12	134.330	30
FS	ID	269-54-13	269-16-56	134.330	30
BS	ID	180-00-27	270-14-47	218.340	10

BS	D	180-00-00	89-45-21	218.340	10
FS	D	269-53-46	90-43-11	134.330	30
FS	ID	89-54-12	269-16-59	134.330	30
BS	ID	0-00-28	270-14-49	218.340	10

Az 20 → 10 = 35°00'00"

JOB: *

TIME: 13:36 DATE: 12-28-1994

Page 2

Side Shot:10-133,Ang-Rt 76.4455,Zenith 90.0600,Slp Dst 96.0400,MW 102
Side Shot:10-134,Ang-Rt 76.4324,Zenith 89.5652,Slp Dst 113.4500,BLD COR J @ EP
HI / HR :Inst H 5.2400,Rod H 9.0000
Side Shot:10-135,Ang-Rt 81.0419,Zenith 88.0520,Slp Dst 111.8100,EP COR
Side Shot:10-136,Ang-Rt 80.3636,Zenith 87.5613,Slp Dst 98.3500,EP COR
HI / HR :Inst H 5.2400,Rod H 7.7000
Side Shot:10-137,Ang-Rt 78.4751,Zenith 88.2213,Slp Dst 73.5100,EP COR
HI / HR :Inst H 5.2400,Rod H 6.7000
Side Shot:10-138,Ang-Rt 88.4115,Zenith 76.4814,Slp Dst 48.9200,EP COR

Easy Survey Raw Data Editor, File ->D:\HAZSITES\ABB\ENRX\TDS\ENRX.RW5

JOB:Name ENRX,Date 11-13-1994,Time 12:49:52
 Mode Setup:North Azimuth,Dist feet,Scale 1.0000,Earth crv OFF
 Store:Pt 1,N 0.0000,E 0.0000,Elv 0.0000,NULL
 Occupy:Occ 20,N 9821.1488,E 9874.7670,Elv 97.1500,PI-20
 Backsight:Occ 20,BS Pt 10,BS azm 35.0000,Back circle 0.0000
 HI / HR :Inst H 5.2400,Rod H 5.2500
 Side Shot:20-100,Ang-Rt 2.5929,Zenith 89.2444,Slp Dst 83.7900,BLD.COR.A/CLF
 Side Shot:20-101,Ang-Rt 80.0436,Zenith 89.2152,Slp Dst 4.1200,BLD COR B
 Side Shot:20-102,Ang-Rt 92.2231,Zenith 90.4041,Slp Dst 37.8700,BLD COR L
 Side Shot:20-103,Ang-Rt 166.1746,Zenith 91.1823,Slp Dst 6.4800,FH. SITE TBM
 HI / HR :Inst H 5.2400,Rod H 0.1800
 Side Shot:20-104,Ang-Rt 232.4414,Zenith 94.4815,Slp Dst 52.0600,MW. 103
 Occupy:Occ 30,N 9744.3065,E 9984.9350,Elv 95.7800,PI-30
 Backsight:Occ 30,BS Pt 20,BS azm 304.5345,Back circle 0.0000
 HI / HR :Inst H 5.0600,Rod H 5.2500
 Side Shot:30-105,Ang-Rt 354.3354,Zenith 85.2745,Slp Dst 18.8400,BLD COR M
 Side Shot:30-106,Ang-Rt 1.0506,Zenith 85.2739,Slp Dst 17.1700,BLD COR C
 Side Shot:30-107,Ang-Rt 101.5426,Zenith 89.3035,Slp Dst 57.3700,BLD COR D
 Side Shot:30-108,Ang-Rt 358.4935,Zenith 89.1838,Slp Dst 129.9000,BLD COR N
 Occupy:Occ 40,N 9849.2551,E 10183.3134,Elv 97.1950,PI-40
 Backsight:Occ 40,BS Pt 30,BS azm 242.0711,Back circle 0.0000
 HI / HR :Inst H 5.0600,Rod H 5.2500
 Side Shot:40-109,Ang-Rt 14.4910,Zenith 89.5311,Slp Dst 135.7700,BLD COR E @ EP
 Side Shot:40-110,Ang-Rt 35.3002,Zenith 89.3141,Slp Dst 130.7200,BLD COR @ EP
 Side Shot:40-111,Ang-Rt 45.1258,Zenith 89.3352,Slp Dst 197.0700,BLD COR H @ EP
 Side Shot:40-112,Ang-Rt 58.5413,Zenith 89.3852,Slp Dst 230.9500,EP COR
 Side Shot:40-113,Ang-Rt 58.1906,Zenith 89.3052,Slp Dst 189.4200,EP COR
 Side Shot:40-114,Ang-Rt 66.2215,Zenith 89.2224,Slp Dst 188.2700,EP COR @ CLF
 Side Shot:40-115,Ang-Rt 66.2529,Zenith 89.3704,Slp Dst 104.7600,PAV CLF ANG.PO
 Side Shot:40-116,Ang-Rt 47.3641,Zenith 88.5059,Slp Dst 22.3900,PAV CLF POST CO
 Side Shot:40-117,Ang-Rt 13.0619,Zenith 89.4134,Slp Dst 68.3100,CLF POST END
 Side Shot:40-118,Ang-Rt 10.1630,Zenith 89.4045,Slp Dst 88.7200,CLF BEG
 Side Shot:40-119,Ang-Rt 7.2317,Zenith 89.5622,Slp Dst 117.1200,CLF END
 Side Shot:40-120,Ang-Rt 3.2132,Zenith 90.0043,Slp Dst 130.5900,EP
 Side Shot:40-121,Ang-Rt 3.4610,Zenith 89.5819,Slp Dst 119.1000,EP
 Side Shot:40-122,Ang-Rt 8.3320,Zenith 89.5510,Slp Dst 119.0400,EP
 Side Shot:40-123,Ang-Rt 20.4113,Zenith 89.3328,Slp Dst 53.2800,EP
 Side Shot:40-124,Ang-Rt 10.1747,Zenith 89.3640,Slp Dst 43.9700,EP
 Side Shot:40-125,Ang-Rt 52.5105,Zenith 88.5508,Slp Dst 33.0400,MW 101
 Side Shot:40-126,Ang-Rt 61.1022,Zenith 89.3818,Slp Dst 197.2500,SS 101
 Occupy:Occ 10,N 10000.0000,E 10000.0000,Elv 98.1000,PI-10
 Backsight:Occ 10,BS Pt 40,BS azm 129.2554,Back circle 0.0000
 HI / HR :Inst H 0.0000,Rod H 0.0000
 Side Shot:10-127,Ang-Rt 85.3207,Zenith 90.0708,Slp Dst 96.8100,EP @ STREET
 Side Shot:10-128,Ang-Rt 85.3126,Zenith 90.0627,Slp Dst 77.8400,EP @ STREET
 Side Shot:10-129,Ang-Rt 85.3421,Zenith 89.4938,Slp Dst 48.0900,EP @ STREET
 Side Shot:10-130,Ang-Rt 86.1117,Zenith 89.4409,Slp Dst 34.6000,EP @ STREET
 Side Shot:10-131,Ang-Rt 8.3556,Zenith 90.0000,Slp Dst 5.2800,CLF END/JOIN#100
 HI / HR :Inst H 5.2400,Rod H 5.2500
 Side Shot:10-132,Ang-Rt 57.3937,Zenith 90.0619,Slp Dst 99.8100,SS 102

Easy Survey Raw Data Editor, File ->D:\HAZSITES\ABB\ENRX\TDS\TRAV.RW5

JOB:Name TRAV,Date 12-28-1994,Time 09:22:21

Mode Setup:North Azimuth,Dist feet,Scale 1.0000,Earth crv OFF

Store:Pt 1,N 0.0000,E 0.0000,Elv 0.0000,NULL

Store:Pt 10,N 10000.0000,E 10000.0000,Elv 98.1000,PI-10

Occupy:Occ 10,N 10000.0000,E 10000.0000,Elv 98.1000,PI-10

Backsight:Occ 10,BS Pt 0,BS azm 215.0000,Back circle 0.0000

Traverse:10-20,Ang-Rt 0.0000,Zenith 90.0000,Slp Dst 218.3370,PI-20

Traverse:20-30,Ang-Rt 89.5345,Zenith 90.4307,Slp Dst 134.3300,PI-30

Traverse:30-40,Ang-Rt 117.1326,Zenith 89.4803,Slp Dst 224.4300,PI-40

Traverse:40-41,Ang-Rt 67.1839,Zenith 89.3926,Slp Dst 237.3500,PI-10

Traverse:41-21,Ang-Rt 85.5345,Zenith 90.1921,Slp Dst 218.3350,PI-20

ATTACHMENT E

FIELD NOTES

Easy Survey Coordinate Editor, File ->ENRX.CR5

Point	Northing	Easting	Elevation	- Description -
1	0.0000	0.0000	0.0000	NULL
10	10000.0000	10000.0000	98.1000	PI-10
20	9821.1488	9874.7670	97.1500	PI-20
30	9744.3065	9984.9350	95.7800	PI-30
40	9849.2551	10183.3134	96.7000	PI-40
100	9887.1805	9926.3406	97.9996	BLD.COR.A/CLF
101	9819.4027	9878.4984	97.1857	BLD COR B
102	9798.1621	9904.8593	96.6918	BLD COR L
103	9815.1129	9872.4142	96.9923	FH. SITE TBM
104	9819.1006	9822.9303	97.8500	MW. 103
105	9753.5435	9968.5826	97.0805	BLD COR M
106	9754.3625	9971.0844	96.9488	BLD COR C
107	9783.5753	10026.7565	96.0809	BLD COR D
108	9816.4174	9876.8999	97.1531	BLD COR N
109	9818.5731	10051.0559	96.7792	BLD COR E @ EP
110	9866.5889	10053.7522	97.5867	BLD COR @ EP
111	9907.9747	9995.2008	98.0081	BLD COR H @ EP
112	9968.2815	9985.4028	97.9297	EP COR
113	9945.2130	10020.0056	98.1152	EP COR
114	9966.4246	10035.9612	98.5691	EP COR @ CLF -
115	9914.5318	10101.3798	97.2089	PAV CLF ANG.POIN
116	9856.8126	10162.2422	96.9595	PAV CLF POST COR
117	9831.8347	10117.2630	96.8763	CLF POST END
118	9822.4215	10098.7501	97.0068	CLF BEG
119	9808.2537	10073.6049	96.6338	CLF END
120	9795.0559	10064.5018	96.4828	EP
121	9800.6024	10074.6041	96.5683	EP
122	9809.8622	10070.9804	96.6774	EP
123	9842.5837	10130.4543	96.9212	EP
124	9835.9720	10141.3988	96.8084	EP
125	9863.2008	10153.3673	97.1334	MW 101
126	9957.5261	10018.4395	97.7551	SS 101
127	9920.6660	9944.5180	97.8991	EP @ STREET
128	9936.2027	9955.4024	97.9540	EP @ STREET
129	9960.6091	9972.4140	98.2450	EP @ STREET
130	9971.8739	9979.8491	98.2595	EP @ STREET
131	9996.0743	10003.5309	98.1000	CLF END/JOIN#100
132	9900.9538	9987.6773	97.9066	SS 102
133	9913.8128	9957.6276	97.9224	MW 102
134	9898.1667	9949.9912	98.1934	BLD COR J @ EP
135	9903.7184	9943.2777	98.0688	EP COR
136	9914.9173	9950.7950	97.8805	EP COR
137	9935.2593	9965.2439	97.7306	EP COR
138	9962.5295	9970.5992	107.8077	EP COR

*Fence Posts
w/NO Fence*

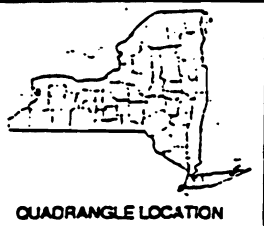
Easy Survey Coordinate Editor, File ->TRAV.CR5

Point	Northing	Easting	Elevation	- Description -
1	0.0000	0.0000	0.0000	NULL
10	10000.0000	10000.0000	98.1000	PI-10
20	9821.1488	9874.7670	97.1500	PI-20
30	9744.3065	9984.9350	95.7800	PI-30
40	9849.2551	10183.3134	96.7000	PI-40

SUMMARY

ENRX INC.				
POINT	NORTHING	EASTING	ELEVATION	DESCRIPTION
SS 101	9957.53	10018.44	97.76	GROUND
SS 102	9900.95	9987.68	97.91	GROUND
MW 101	9863.20	10153.37	N/A	FLUSH MOUNT
			97.13	CASE
			96.51	RISER
MW 102	9913.81	9957.63	N/A	FLUSH MOUNT
			97.93	CASE
			97.71	RISER
MW 103	9819.10	9822.93	N/A	FLUSH MOUNT
			97.87	CASE
			97.33	RISER

ATTACHMENT D
TABULATION OF DATA



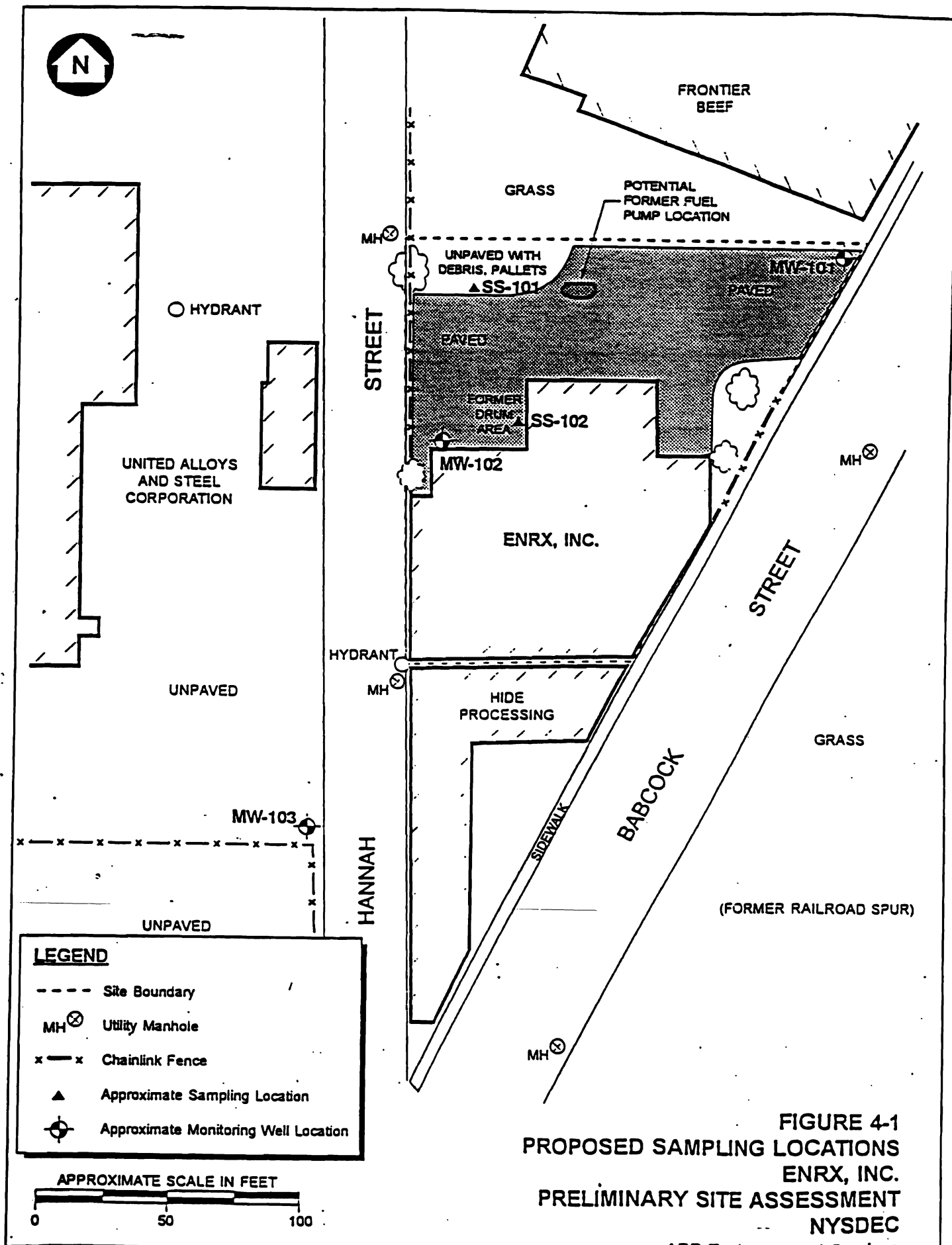
SOURCE: N.Y.S. DEPARTMENT OF TRANSPORTATION 7.5-MINUTE SERIES QUADRANGLES;
BUFFALO NE, NEW YORK, 1989, AND BUFFALO SE, NEW YORK, 1989.



SITE NO.: 915150
LOCATION: CITY OF BUFFALO
ERIE COUNTY



FIGURE 1-1
SITE LOCATION MAP
ENRX, INC.
PRELIMINARY SITE ASSESSMENT
NYSDEC



ENRX INC.

+ π - ELEV.

DATUM

2.445

5.12

4.58

5.295

5.82

4.97

5.76

6.15

5.93

7.35

6.73

7.16

5.275

6.195

6.815

2.57

26.115

26.110

SITE BLVD F11 E/SIDE OF HANNAH ST
BY RP (NYT-39) NORTH B-BOULT

CHISEL "X"

RISE RIM (FLUSH)

RISE

RIM (FLUSH)

RISE

RIM (FLUSH)

PI-30

PI-40

MW-101

MW-102

PI-10

PI-20

MW-103

SITE BM

