

### APPENDIX L

### LEGGETTE, BRASHEARS & GRAHAM 1991



# APPENDIX L1 LBG 1991 SOIL BORING LOGS

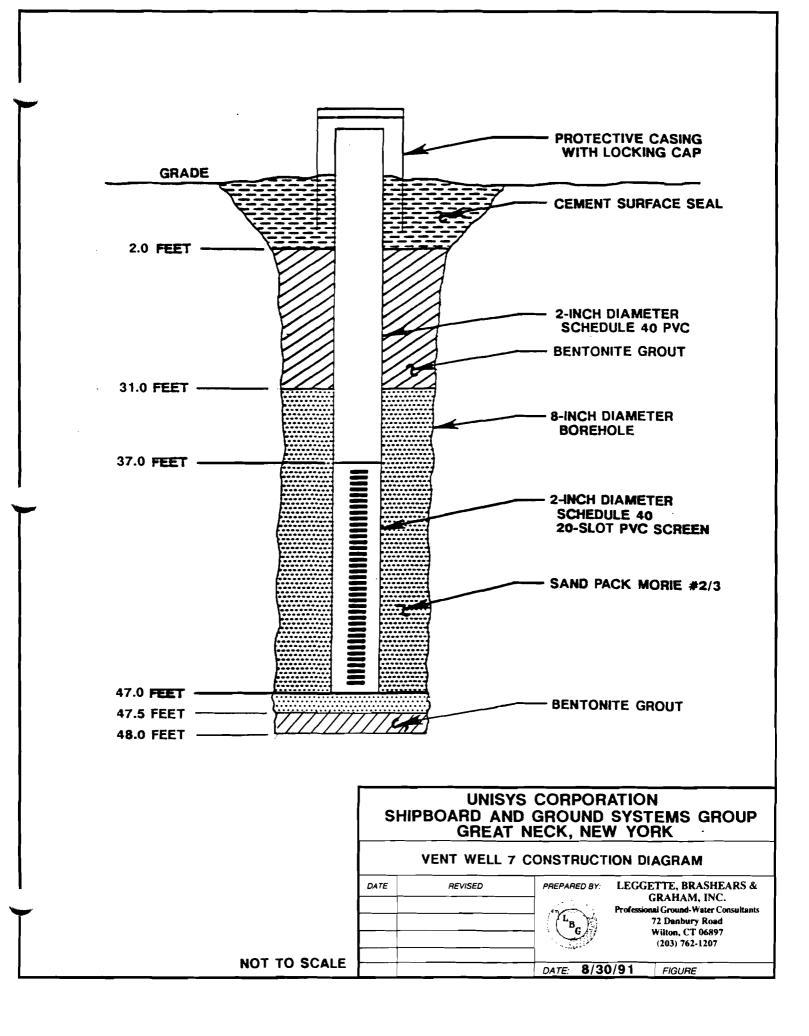
## **GEOLOGIC LOG**

OWNER Unisys Corporation, Great Neck, NY LEGGETTE, BRASHEARS & GRAHAM, INC. | WELL NO. Soil Boring 10/Vent Well 7

	Soil Boring 10/Vent Well /			
WILTON, CONNECTICUT	PAGE 1 OF 1 PAGES			
LOCATION South side of building west of	SCREEN TYPE PVC			
reclamation area	DIAM. 2-inch SLOT NO. 20			
DATE COMPLETED July 30, 1991	SETTING 37 - 47 ft bg			
DRILLING	SAND PACK Morie No. 2/3			
COMPANY R & L Well Drilling	CASING PVC			
DRILLING METHOD Hollow-stem augers				
SAMPLING	SETTING 2.5 ft ag to 37 ft bg			
METHOD Split spoon	DEVELOPMENT			
OBSERVER Keith Yocis	DURATION			
REFERENCE POINT (RP) Grade	STATIC WATER LEVEL			
ELEVATION OF RP	YIELD			

REMARKS

DEPTH (FEE FROM	T)   TO	DESCRIPTION	
10	12	SAND, medium, some coarse, little fine; and gravel, fine to medium;	
		trace gravel, coarse; and silt; brown; PID = 3.6 ppm; 0.9-foot	
		recovery.	
20	22	SAND, medium; some coarse; and gravel, fine; little sand, fine to	
-		very coarse; and gravel, medium to coarse; trace cobble; brown;	
		PID = 29 ppm; 1.3-foot recovery.	
30	32	SAND, medium; some fine to coarse; little gravel, fine to coarse; and	
		cobbles; and sand, very fine to very coarse; tan; PID = 5.5 ppm;	
		1.3-foot recovery.	
40	42	SAND, medium; some coarse; little fine; and gravel, fine to coarse;	
		trace cobble; PID = 5.8 ppm; 0.8-foot recovery.	
1	48	End of Borehole.	



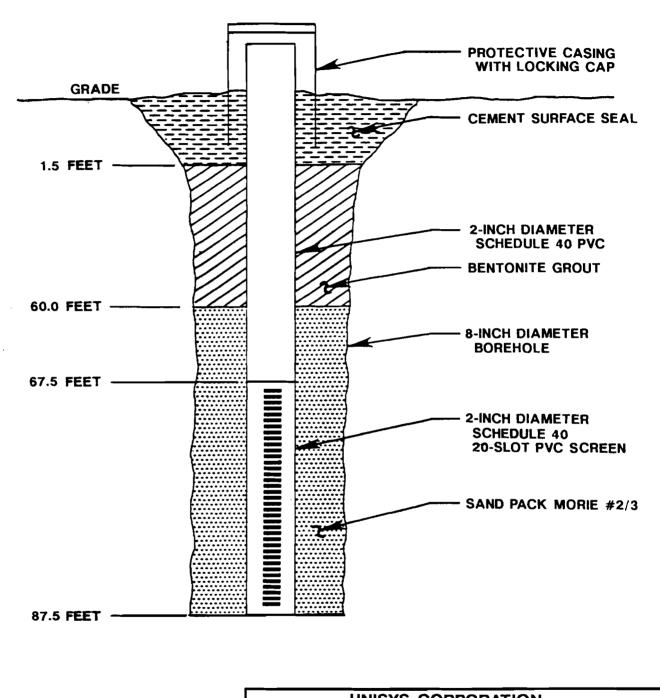
GEOLOGIC LOG	OWNER Unisys Corporation, Great Neck, NY			
LEGGETTE, BRASHEARS & GRAHAM, INC.	WELL NO. Soil Boring 11/Vent Well 8			
WILTON, CONNECTICUT	PAGE 1 OF 2 PAGES			
LOCATION	SCREEN TYPE PVC			
	DIAM. 2-inch SLOT NO. 20			
DATE COMPLETED July 24, 1991	<b>SETTING</b> 67.5 - 87.5 ft bg			
	<b>SAND PACK</b> Morie 2/3 60 - 87.5 ft bg			
<del></del>	CASING PVC			
	accento.			
SAMPLING	SETTING 2.5 ft ag to 67.5 ft bg			
METHOO Split spoon	DEVELOPMENT			
OBSERVER Keith Yocis	DURATION			
REFERENCE POINT (RP) Grade	STATIC WATER LEVEL			
ELEVATION OF RP	YIELD			
METHOD Split spoon	SAND PACK Morie 2/3 60 - 87.5 ft bg  CASING PVC  SETTING 2.5 ft ag to 67.5 ft bg  DEVELOPMENT  DURATION  STATIC WATER LEVEL			

REMARKS

DEPTH (FEET) FROM TO		DESCRIPTION			
10	12	SAND, medium to coarse, some fine, and fine gravel; little medium to			
		coarse gravel; trace cobbles; brown.			
20	22	SAND, fine to medium, little very fine to coarse; some fine gravel,			
		trace medium to coarse; brown, PID = 5.7 ppm, 1.1-foot recovery.			
30	32	SAND, medium, some fine to coarse, little very fine; some fine gravel,			
		trace medium to coarse; some silt; brown; PID = 61.9 ppm; 1.0-foot			
		recovery.			
33	38	CLAY and Sand, fine to medium; brown (cuttings).			
40	42	SAND, fine to medium, little very fine, and silt; trace coarse sand,			
		and cobbles; brown; PID = 10.1 ppm; 0.7-foot recovery.			
	<del></del> -	·			

OWNER	Unisys Corporation, Great Ne	ck, New York				
WELL NO.	Soil Boring 11/Vent Well 8	PAGE	2	OF	2	PAGES

WELL NO		Soring 11/Vent Well 8 PAGE 2 OF 2 PAGES			
DEPTH (F FROM	EET)	DESCRIPTION			
50	52	SAND, fine to medium, some coarse; little fine to coarse gravel;			
		trace silt and clay; gray-brown; 0.5 foot thick.			
		SAND, medium to coarse, some very coarse, trace fine; little fine to			
		coarse gravel; brown; 0.9 foot thick; PID = 11.6 ppm; 1.4-foot			
		recovery.			
60	62	SAND, fine to medium, little very fine to coarse; trace silt and			
		fine gravel; brown; PID = 11.5 ppm; 0.7-foot recovery.			
70	72	SAND, medium to coarse, some very coarse, trace fine; little fine to			
		coarse gravel; brown; 0.7 foot thick.			
		SAND, medium to very coarse, little fine; some fine gravel; tan;			
		0.25 foot thick.			
		GRAVEL, fine; some very coarse to fine sand; iron stained red-orange;			
		0.15 foot thick.			
		SAND, medium to very coarse, little fine; some fine to medium gravel;			
		trace micas; tan; 0.2 foot thick; PID = 5.2 ppm; 1.3-foot recovery.			
80	82	SAND, medium to very coarse, little fine; some fine to coarse gravel;			
		trace micas; saturated; tan; PID = 189 ppm; 1.6-foot recovery.			
	90	End of Borehole.			
	1				



### **VENT WELL 8 CONSTRUCTION DIAGRAM**

DATE REVISED PREPARED BY:

**NOT TO SCALE** 

LEGGETTE, BRASHEARS & GRAHAM, INC.

Professional Ground-Water Consultants 72 Danbury Road Wilton, CT 06897 (203) 762-1207

DATE: 8/20/91

FIGURE

### **GEOLOGIC LOG** OWNER Unisys Corporation, Great Neck, NY LEGGETTE, BRASHEARS & GRAHAM, INC. WELL NO. Soil Boring 12 WILTON, CONNECTICUT PAGE OF 1 PAGES 1 SCREEN TYPE LOCATION West of the Reclamation area north of the garage DIAM. SLOT NO. DATE COMPLETED July 25, 1991 SETTING SAND PACK DRILLING COMPANY R & L Well Drilling CASING DRILLING METHOD Hollow-stem auger SETTING SAMPLING Split spoon METHOD DEVELOPMENT OBSERVER Keith Yocis DURATION Grade REFERENCE POINT (RP) STATIC WATER LEVEL ELEVATION OF RP YIELD

REMARKS

No well installed.

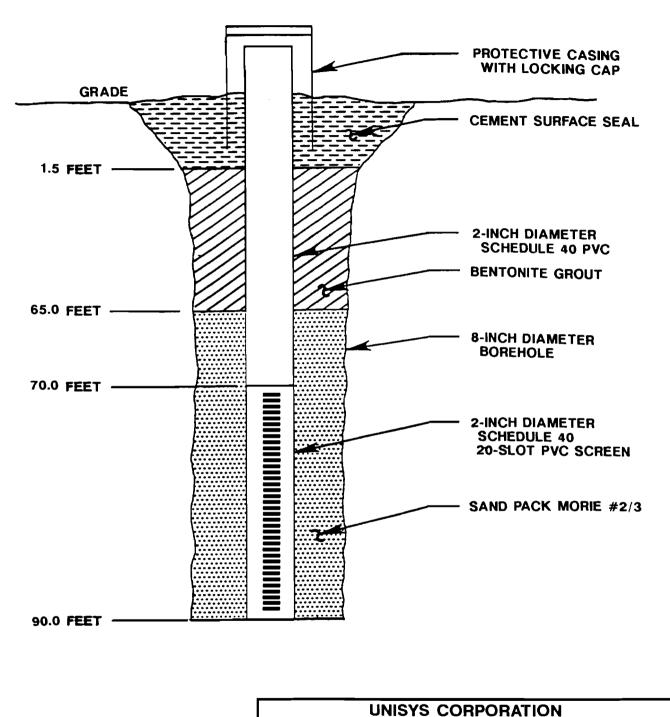
<u> </u>			
DEPTH (FEET	л) то	DESCRIPTION	
0	2.5	Hit refusal (boulder) move hole.	
0	3	Hit refusal (boulder) move hole.	
0	2	Hit refusal, abandon hole.	
		<u> </u>	

GEOLOGIC LOG LEGGETTE, BRASHEARS & GRAHAM, INC.	OWNER Unisys Corporation, Great Neck, NY WELL NO. Soil Boring 13/Vent Well 9				
WILTON, CONNECTICUT	PAGE 1 OF 2 PAGES				
LOCATION	SCREEN TYPE PVC				
	DIAM. 2-inch SLOT NO. 20				
DATE COMPLETED July 26, 1991	SETTING 70 - 90 ft bg				
DRILLING	SAND PACK Morie 2/3 65 - 90 ft bg				
COMPANY R & L Well Drilling	- CASING PVC				
DRILLING METHOD Hollow-stem augers	SETTING OF G				
SAMPLING	SETTING 2.5 ft ag to 70 ft bg				
METHOO Split spoon	DEVELOPMENT				
OBSERVER Keith Yocis	DURATION				
REFERENCE POINT (RP) Grade	STATIC WATER LEVEL				
ELEVATION OF RP	YIELO				
REMARKS					

r			
DEPTH (FEET	n)   то	DESCRIPTION	
10	12	SAND, medium to very coarse, little fine; trace fine gravel; tan;	
	_	0.5 foot thick.	
		SAND, medium to coarse, some fine, trace very coarse; some fine gravel;	
		tan with brown bands; 0.4 foot thick; PID = 26.3 ppm; 0.9-foot	
		recovery.	
30	32	Clay and sand, fine to medium; gradational to all clay; brown; 0.8	
		foot thick.	
		SAND, medium; trace fine gravel; brown; 0.1 foot thick.	
		CLAY; little fine to medium sand; trace micas; brown; 0.4 foot thick.	
	_	SAND, fine to coarse; tan; 0.1 foot thick.	
		CLAY; trace fine sand; brown; 0.1 foot thick.	
		SAND, medium, little coarse to fine; tan; 0.3 foot thick;	

OWNER	Unisys Corporation, Great Neck, New York	_			
WELL NO.	Soil Boring 13/Vent Well 9	PAGE 2	OF	2	PAGES

WELL NO. Soil Boring 13/Vene Well 9 PAGE 2 OF 2 PAGES					
DEPTH (FEET) FROM TO DESCRIPTION					
50	52	SAND, medium, some fine to coarse; trace fine gravel and cobbles;			
		tan; PID = 10.9 ppm; 0.7-foot recovery.			
60	62	SAND, medium, some fine to coarse, trace very coarse; some fine			
		gravel; tan; PID = 10.9 ppm; 1.0-foot recovery.			
70	72	SAND, medium to coarse, some very coarse; little fine to coarse			
		gravel; some iron staining; tan; 0.8 foot thick.			
		SAND, medium, some fine, little very coarse; trace fine gravel; tan;			
		0.7 foot thick; PID = 9.0 ppm; 1.5-foot recovery.			
80	82	SAND, medium to very coarse, some fine; little fine to medium gravel;			
		trace micas; saturated; tan; PID = 29.5 ppm; 1.5-foot recovery.			
	90	End of Borehole.			



### **VENT WELL 9 CONSTRUCTION DIAGRAM**

DATE REVISED PREPARED BY:

LEGGETTE, BRASHEARS & GRAHAM, INC.
Professional Ground-Water Consultants

rofessional Ground-Water Consultan 72 Danbury Road Wilton, CT 06897 (203) 762-1207

DATE: 8/20/91

FIGURE

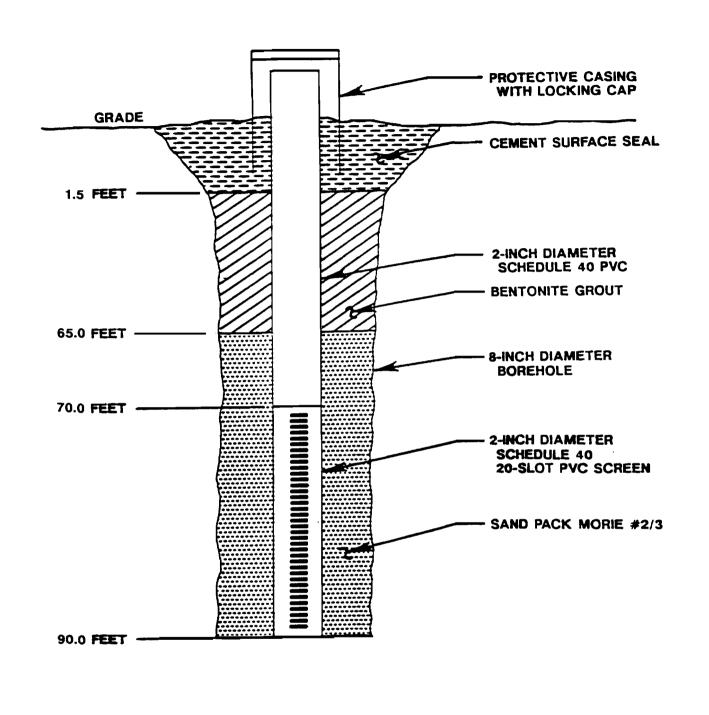
GEOLOGIC LOG LEGGETTE, BRASHEARS & GRAHAM, INC.	OWNER Unisys Corporation, Great Neck, NY WELL NO. Soil Boring 14/Vent Well 10		
WILTON, CONNECTICUT	PAGE 1 OF 2 PAGES		
LOCATION Approximately 190 feet north of	SCREEN TYPE PVC		
reclamation area, 19 feet east of building	DIAM. 2-inch SLOT NO. 20		
DATE COMPLETED August 16, 1991	SETTING 70 - 90 ft bg		
DRILLING COMPANY R & L Well Drilling	SAND PACK Morie No. 2/3 65 - 90 ft bg		
DRILLING METHOD Hollow-stem augers	CASING PVC  SETTING 2.5 ft ag to 65 ft bg		
SAMPLING METHOD Split spoon - 3 inch	DEVELOPMENT		
OBSERVER Robert Emig	DURATION		
REFERENCE POINT (RP) Grade	STATIC WATER LEVEL		
ELEVATION OF RP	YIELD		

### REMARKS

DEPTH (FEET	TO TO	DESCRIPTION
0	20	Sand, fine, and silt; large cobbles; brown (cuttings).
20	22	SAND, fine to medium; some medium gravel and silt; brown; 0.8 foot thick.
		SAND, medium; trace silt; yellow-brown; 0.5 foot thick; PID = 7.1 ppm.
22	30	SAND, fine to medium; some medium gravel; little silt; dark brown
		(cuttings).
30	32	Clay and Sand, medium; little medium gravel; brown; 0.6-foot recovery;
		PID = 4.8 ppm.
32	40	Sand, medium, and clay; some medium gravel; brown (cuttings).
40	42	Sand, medium, and clay; little medium gravel; brown; 1.4-foot recovery;
		PID = 4.6 ppm.
50	52	SAND, medium; some silt and coarse gravel; brown-yellow; 1.4-foot
		recovery; PID = 3.9 ppm.

OWNER	Unisys Corporation, Great Neck, New York					_
WELL NO.	Soil Boring 14/Vent Well 10	PAGE	2	OF	2	PAGES

WELL NO.	Soil B	oring 14/Vent Well 10	PAGE 2	OF 2 PAGE	:s
DEPTH (FE	≣ΕΤ)   ΤΟ	DESCRI	PTION		
52	60	SAND, coarse; some silt; wet at appro	oximately 55 feet	(cuttings).	
60	62	SAND, medium to coarse; some coarse	gravel; trace sil	t; yellow-brown;	
		1.9-foot recovery; PID = 4.0 ppm	•		
62	70	Sand, coarse and gravel, coarse; some	e silt; brown (cu	ttings).	
<b>7</b> 0	72	SAND, medium; little fine gravel and	silt; brown-yell	ow; 1.4 feet	
		thick.			
		SAND, coarse and gravel, coarse; trac	ce silt; tan; 0.6	foot thick;	
		PID = 5.9 ppm.			
72	80	Sand, coarse and gravel, medium; some	e silt: brown (cu	ttings).	
80	82	SAND, medium to coarse; little grave	l; trace silt; ta	n-brown mix;	
	ļ	saturated; 2-foot recovery; PID =	= 5.2 ppm.		
80	90	SAND, coarse and gravel, medium; trac	ce silt (cuttings	).	
	90	End of Borehole.			
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### **VENT WELL 10 CONSTRUCTION DIAGRAM**

DATE REVISED PREPARED BY

LEGGETTE, BRASHEARS & GRAHAM, INC.

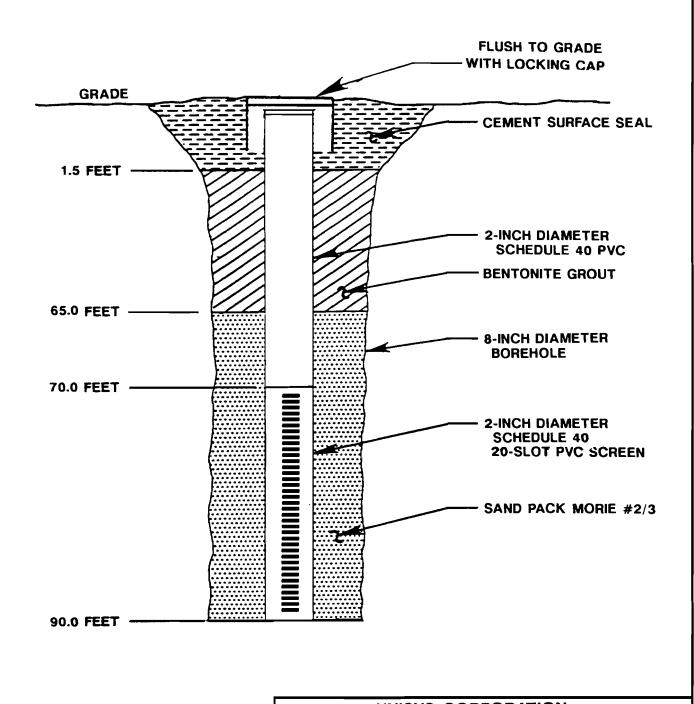
Professional Ground-Water Consultants 72 Danbury Road Wilton, CT 06897 (203) 762-1207

DATE: 8/27/91 FIGURE

GEOLOGIC LOG LEGGETTE, BRASHEARS & GRAHAM, INC. WILTON, CONNECTICUT	OWNER Unisys Corporation, Great Neck, NY WELL NO. Soil Boring 15/Vent Well 11 PAGE 1 OF 2 PAGES
LOCATION South parking lot, east of	SCREEN TYPE PVC
Guard Post 3	DIAM. 2-inch SLOT NO. 20
DATE COMPLETED August 6, 1991	SETTING 70 - 90 ft bg
DRILLING COMPANY R & L Well Drilling DRILLING	SAND PACK Morie 2/3 65 - 90 ft bg  CASING PVC
METHOD Hollow-stem augers	SETTING 0 to 70 ft bg
SAMPLING METHOD Split spoon	DEVELOPMENT
OBSERVER John Benvegna	DURATION
REFERENCE POINT (RP) Grade	STATIC WATER LEVEL
ELEVATION OF RP	YIELD
REMARKS	<del></del>

DEPTH (FEE FROM	г)   то	DESCRIPTION
0	0.3	Asphalt.
0.3	10	SAND, coarse to very coarse and Gravel; some cobbles and boulders;
		brown (cuttings).
10	12	Sand, coarse to very coarse, little medium, and fine to medium gravel,
		trace coarse; brown-tan; PID = 2.4 ppm; 1.4-foot recovery.
12	20	SAND, coarse to very coarse and gravel; brown (cuttings).
20	22	Sand, coarse to very coarse, little medium, and fine to coarse gravel;
		brown; PID = 1.7 ppm; 1.8-foot recovery.
30	32	SAND, medium to coarse, little fine; little fine to medium Gravel;
		brown; PID = 4.2 ppm; 2.0-foot recovery.

	OWNER	Unisys	Corporation, Great Neck, New York		
	WELL NO. Soil Boring 15/Vent Well 11 PAGE 2 OF 2 PAGES				
DEPTH (FEET) FROM TO DESCRIPTION			DESCRIPTION		
	40	42	Sand, medium, some fine, little coarse to very coarse, and fine		
			gravel; trace medium to coarse gravel; brown; PID = 7.8 ppm;		
			1.3-foot recovery.		
	50	52	SAND, medium, some coarse, little very coarse to fine; trace fine		
$\int$			gravel; brown; PID = 3.0 ppm; 0.3-foot recovery.		
	60	62	SAND, medium, some coarse, little very coarse; some fine to medium		
			gravel; brown; PID = 9.5 ppm; 1.4-foot recovery.		
	70	72	SAND, medium to very coarse; some fine to medium gravel; trace fine		
			sand and coarse gravel; iron stains; brown; PID = 4.1 ppm;		
			0.8-foot recovery.		
	80	82	SAND, very coarse, little medium, trace fine; some fine to medium		
			gravel, trace coarse; saturated; brown; PID = 7.8 ppm; 1.4-foot		
			recovery.		
		90	End of Borehole.		
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### **VENT WELL 11 CONSTRUCTION DIAGRAM**

DATE REVISED PREPARED BY.

LEGGETTE, BRASHEARS & GRAHAM, INC.
Professional Ground-Water Consultants

72 Danbury Road Wilton, CT 06897 (203) 762-1207

DATE: 8/20/91

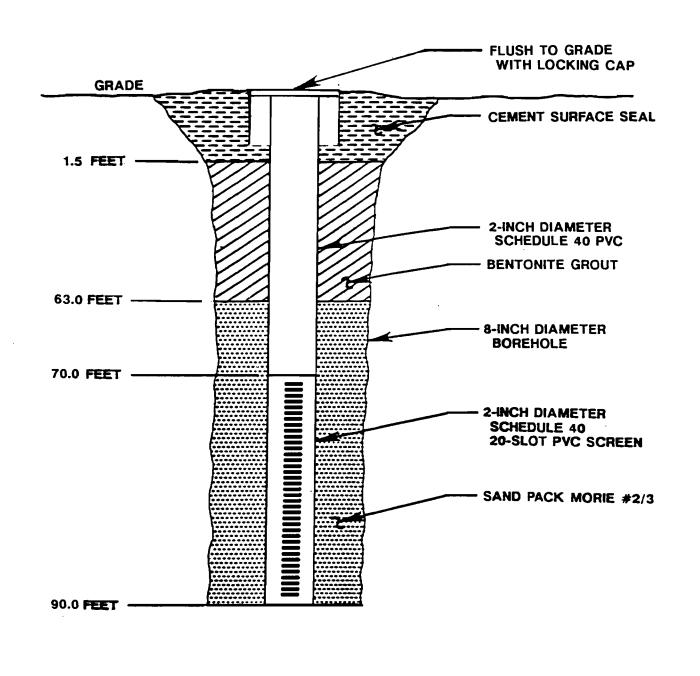
FIGURE

GEOLOGIC LOG LEGGETTE, BRASHEARS & GRAHAM, INC. WILTON, CONNECTICUT	OWNER Unisys Corporation, Great Neck, NY  WELL NO. Soil Boring 16/Vent Well 12  PAGE 1 OF 2 PAGES
LOCATION South of Guard Post 3, 6 ft north	SCREEN TYPE PVC
of culvert.	DIAM. 2-inch SLOT NO. 20
DATE COMPLETED August 14, 1991	SETTING 70 - 90 ft bg
DRILLING COMPANY R & L Well Drilling	SAND PACK Morie No. 2; 63 - 90 ft bg
DRILLING METHOD Hollow-stem augers	CASING PVC
SAMPLING METHOD Split spoon - 3 inch	DEVELOPMENT
OBSERVER Robert Emig	DURATION
REFERENCE POINT (RP) Grade	STATIC WATER LEVEL
ELEVATION OF RP	YIELD
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DEPTH (FEET)		DECORIDATION
FROM	ТО	DESCRIPTION
0	20	SAND, medium; little silt; cobbles; dark brown (cuttings).
20	22	SAND, medium; some silt; little medium gravel; dark brown; 1.4-foot
		recovery; PID = 33.5 ppm.
22	30	SAND, medium; some very coarse gravel; trace silt; brown (cuttings).
30	32	No recovery.
32	40	SAND, medium; little silt; gravel; brown (cuttings).
40	42	SAND, medium; some silt; little coarse gravel; dark brown; 0.6 foot thick
	-	SAND; coarse; trace coarse gravel; yellow-brown; 0.9 foot thick;
		PID = 39.8 ppm.
42	50	SAND, medium; trace silt and gravel; dark brown (cuttings).
50	52	SAND, fine to medium; little fine gravel; yellowish; 1.4-foot
		recovery; PID = 14.1 ppm.

OWNER	Unisys	Corporation, Great Neck, New Yor	k			
WELL NO.	Soil Bo	oring 16/Vent Well 12	PAGE 2	OF 2	PAGES	
DEPTH (FE	ET) TO	DESCRIPTION				
52	60	SAND, medium; trace silt; brown	(cuttings).			
60	62	SAND, fine to medium; trace sil	t; brownish-yellow;	moist;		
		1.3 feet thick; PID 26.4 pp	m.			
	T	SAND, fine; some silt; brown; m	oist; 0.7 foot thick	ζ.		
70	72	SAND, medium; some coarse grave	l; trace silt yellow	wish-brown;		
		2-foot recovery; PID = 21.8	ppm.			
72	80	SAND, medium; some medium grave	l; trace silt; light	t brown (cutt:	ings).	
80	82	SAND, medium, and coarse gravel	; trace silt; brown-	-rust; satura	ted;	
		1.5-foot recovery; PID = 32	.8 ppm.			
82	90	SAND, medium, and medium gravel	; some silt; brown	(cuttings).		
	90	End of Borehole.				



### **VENT WELL 12 CONSTRUCTION DIAGRAM**

DATE REVISED PREPARED BY:

LEGGETTE, BRASHEARS & GRAHAM, INC.
Professional Ground-Water Consultants
72 Danbury Road
Wilton, CT 06897

Wilton, CT 06897 (203) 762-1207

DATE: 8/27/91 FIGURE

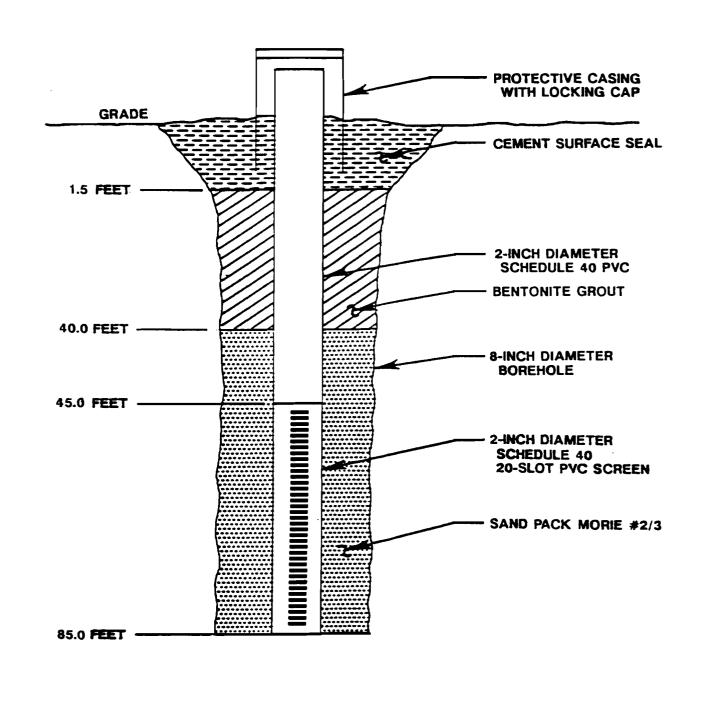
#### **GEOLOGIC LOG** OWNER Unisys Corporation, Great Neck, NY LEGGETTE, BRASHEARS & GRAHAM, INC. WELL NO. Soil Boring 17/Vent Well 13 WILTON, CONNECTICUT PAGE PAGES SCREEN TYPE LOCATION PVC Reclamation Area DIAM. SLOT NO. 2-inch 20 DATE COMPLETED August 22, 1991 SETTING 45 - 85 ft bg SAND PACK DRILLING Morie No. 2; 40 - 90 ft bg COMPANY R & L Well Drilling CASING PVC DRILLING METHOD Hollow-stem augers SETTING 2.5 ft ag to 45 ft bg SAMPLING METHOO split spoon - 3 inch DEVELOPMENT OBSERVER Robert Emig DURATION REFERENCE POINT (RP) Grade STATIC WATER LEVEL ELEVATION OF RP YIELD

REMARKS

DEPTH (FEE	л)   то	DESCRIPTION
0	10	SAND, medium to coarse; some coarse gravel; light brown (cuttings).
10	12	Sand, medium, and gravel, coarse; some silt; brown; 0.4-foot recovery.
12	20	Sand, medium, and gravel, coarse; some cobbles; trace silt; brown
	_	(cuttings).
20	22	Sand, medium, and gravel, coarse; some silt; brown; 2-foot recovery;
		PID = 4,689 ppm.
22	30	SAND, medium; some silt; little medium gravel; brown (cuttings);
30	32	SAND, medium; some silt; little gravel; 1.2 feet thick.
	_	Clay and Sand, fine to medium; little medium gravel; brown; 0.8 foot
		thick; PID = 5,300 ppm.
32	40	Sand, medium, and gravel, very coarse; trace silt; brown (cuttings).

OWNER	Unisys	Corporation, Great Neck, New York		
WELL NO. Soil Boring 17/Vent Well 13 PAGE 2 OF 2 PAGE				
DEPTH (FE FROM	ET)   TO	DESCRIPTION		
40	42	Sand, medium, and clay; some coarse gravel; brown; 2-foot recovery;		
		PID = 1,307 ppm.		
42	45	SAND, medium; some silt; brown (cuttings).		
45	50	SAND, medium; some silt; gray; heavy odor (cuttings).		
50	52	SAND, medium; little coarse gravel; gray; heavy odor; 1.4 feet thick.		
		Sand, medium, and silt; some clay; little coarse gravel; brown; heavy		
		odor; 0.6 foot thick; PID = 4,617 ppm.		
52	60	SAND, medium; some clay; little gravel; trace silt; gray-brown		
		(cuttings).		
60	62	SAND, medium to coarse; little fine gravel and silt; brownish-gray;		
<b></b>	<u> </u>	odor; 2-foot recovery; PID = 4,200 ppm.		
62	70	SAND, medium to coarse; trace medium gravel and silt; gray (cuttings).		
70	72	SAND, medium to coarse; some coarse gravel and silt; dark and light		
		gray; 2-foot recovery; PID = 3,620 ppm.		
72	80	SAND, medium to coarse; little coarse gravel; trace silt; gray		
<u> </u>		(cuttings).		
80	82	SAND, medium to coarse; little medium gravel; trace silt; black-		
		gray; saturated; odor; 1.6-foot recovery; PID 2,500 ppm.		
82	90	SAND, coarse; some coarse gravel; little silt; gray (cuttings).		
	90	End of Borehole.		
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### **VENT WELL 13 CONSTRUCTION DIAGRAM**

DATE REVISED PREPAREO BY: LEGGETTE, BRASHEARS & GRAHAM, INC.

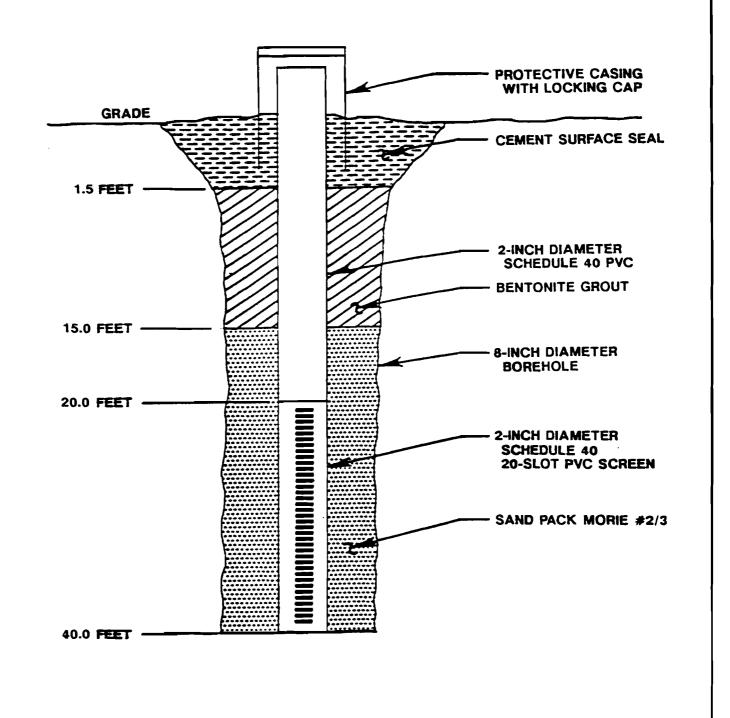
Professional Ground-Water Consultants 72 Danbury Road Wilton, CT 06897 (203) 762-1207

DATE: 8/27/91 FIGURE

### **GEOLOGIC LOG** OWNER Unisys Corporation, Great Neck, NY LEGGETTE, BRASHEARS & GRAHAM, INC. WELL NO. Soil Boring 18/Vent Well 14 OF 1 WILTON, CONNECTICUT PAGE **PAGES** LOCATION Reclamation Area SCREEN TYPE PVC DIAM. SLOT NO. 2-inch 20 SETTING 20 - 40 ft bg DATE COMPLETED August 23, 1991 SAND PACK Morie No. 2; 15 - 40 ft bg DRILLING COMPANY R & L Well Drilling CASING PVC DRILLING METHOD Hollow-stem augers SETTING 2.5 ft ag to 20 ft bg SAMPLING METHOD Grab/cuttings DEVELOPMENT DURATION OBSERVER Robert Emig REFERENCE POINT (RP) Grade STATIC WATER LEVEL ELEVATION OF RP YIELD

### REMARKS

DEPTH (FEET) FROM TO		DESCRIPTION					
0	10	SAND, medium; some coarse gravel and large cobbles; trace silt;					
	_	brown.					
10	15	SAND, medium; some coarse gravel; large cobbles; trace silt; brown.					
15	20	SAND, medium; some medium gravel; trace silt; brown; odor.					
20	30	SAND, medium; little medium gravel; trace silt; brown, odor; PID =					
		4,400 ppm.					
30	40	SAND, medium; little coarse gravel; trace silt; dark-brown; heavy					
		odor.					
40	42	Sand, coarse, and gravel; some silt; gray; heavy odor, 0.3 foot thick.					
-		Clay and sand, medium; brown; saturated; heavy odor; 1.7-foot thick;					
		PID = 4,100 ppm. (split spoon).					
	42	End of Borehole.					



### **VENT WELL 14 CONSTRUCTION DIAGRAM**

DATE REVISED PREPARED BY:

LEGGETTE, BRASHEARS & GRAHAM, INC.

Professional Ground-Water Consultants 72 Danbury Road Wilton, CT 06897 (203) 762-1207

DATE: 8/27/91

FIGURE

# APPENDIX L2 LBG 1991 SOIL CHEMISTRY



### an environmental testing company-

200 Monroe Turnpike Monroe, Connecticut 06468 (203) 261-4458 FAX (203) 268-5346

### REPORT TRANSMITTAL

REPORT NUMBER 30910-1527

DATE August 22, 1991

CLIENT

Unisys Corporation 3199 Pilot Knob MS F1B05 Eagan, MN 55121

ATTENTION \_\_\_ Mr. Kevin Krueger

The above referenced report is enclosed. Copies of this report and supporting data will be retained in our files in the event they are required for future reference.

If there are any questions concerning this report, please do not hesitate to contact us.

Any samples submitted to our Laboratory will be retained for a maximum of sixty (60) days from receipt of this report, unless other arrangements are desired.

30910-1527 UNISYS CORPORATION 3199 Pilot Knob MS F1B05 Eagan, Minnesota 55121

Re: Great Neck, New York

Attention: Mr. Kevin Krueger

### **PURPOSE**

Ten samples collected on July 23, 24 and 26, 1991 were submitted to IEA, Inc. by Unisys Corporation. The client requested the samples be analyzed for TCL volatile organics plus a library search for non-target compounds.

### METHODOLOGY

Volatile organics were determined using purge and trap GC/MS. The instrumentation used was a Tekmar Dynamic Headspace Concentrator interfaced with a Hewlett-Packard Model 5995C GC/MS/DS.

All analyses were conducted according to NYSDEC '89 Protocols.

### DISCUSSION

<u>Volatile Organics</u> - The laboratory followed the USEPA CLP SOW Document #OLMO1.0 for the GC/MS calibration criteria.

Due to instrumentation problems, samples SVB112022, SVB114042, SVB116062, SVB118082 and TB 07/25/91 were not run within the seven day holding time. The client and NYSDEC were contacted regarding the holding time issue and permission was granted to proceed with the analyses. The samples were run on 08/03/91 (out of holding time) and these runs were reported. However, sample TB 07/25/91 leaked during the 08/03/91 run and could not be rerun due to insufficient sample.

Batch QC has been provided.

### RESULTS

The results are presented in the following Tables. Also enclosed are the data packages containing all relevant QA/QC and raw data.

Jahone

Prepared by:

Laboratory Manager

JCC/mt

cc: R. Vitale (Environmental Standards)

The liability of IEA, Inc. is limited to the actual dollar value of this project.

### TABLE 1.0 30910-1527 UNISYS CORPORATION **EPA TCL VOLATILE ORGANICS**

All values are ug/L.

Sample Identification

<u>Dilution Factor</u> <u>1.0</u> <u>1.0</u>	
Method Blank I.D. >B4946 >B4946	
CompoundMethodTBBlank07/27/91	Quantitation Limits with no <u>Dilution</u>
	Dilution  10 10 10 10 10 10 5 10 5 5 5 5 5 5 5 5
Styrene U U Xylene (total) U U	5 5

U - See Appendix for definition. Note: Sample detection limit = quantitation limit x dilution factor.

### TABLE 1.1 30910-1527 UNISYS CORPORATION **EPA TCL VOLATILE ORGANICS**

All values are ug/Kg.

### Sample Identification

<u>Dilution Factor</u>	1.00	1.02	1.11	1.04	1.09	
Method Blank I.D.	<u>&gt;B4962</u>	<u>&gt;B4962</u>	>B4962	<u>&gt;B4962</u>	<u>&gt;B4962</u>	
<u>Compound</u>	Method <u>Blank</u>		SVB 114042	SVB 116062	SVB 118082	Quantitation Limits with no <u>Dilution</u>
Chloromethane Bromomethane Vinyl Chloride Chloroethane Methylene Chloride Acetone Carbon Disulfide 1,1-Dichloroethene 1,1-Dichloroethane 1,2-Dichloroethene (total)  hloroform ,2-Dichloroethane 2-Butanone 1,1,1-Trichloroethane Carbon Tetrachloride	0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0	טטטטטטטטטטטטטטטט	000000001 0000001	טטטטטטטטטטטטטטטט	V U U U U U U U U U U U U U U U U U U U	10 10 10 10 5 10 5 5 5 5 5 5 5
Vinyl Acetate Bromodichloromethane 1,2-Dichloropropane cis-1,3-Dichloropropene Trichloroethene Dibromochloromethane	U U U U U	U U U U U	บ บ บ บ บ <b>9</b>	V U U U U	U U U U	10 5 5 5 5 5
1,1,2-Trichloroethane Benzene trans-1,3-Dichloropropene Bromoform 4-Methyl-2-pentanone 2-Hexanone Tetrachloroethene 1,1,2,2-Tetrachloroethane Toluene Chlorobenzene Ethylbenzene Styrene	טטטטטטטטטטטטטטטט	טטטטטטטטטטטטטטטטטטטטטטטטטטטטטטטטטטטטטטט	0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0	טטטטטטטטטטטטטטטט		5 5 5 5 10 10 5 5 5 5 5
Xylene (total)	Ŭ	Ŭ	Ŭ	Ŭ	ŭ	5

U, J, B - See Appendix for definition.
'ote: Sample detection limit = quantitation limit x dilution factor.

### TABLE 1.2 30910-1527 UNISYS CORPORATION EPA TCL VOLATILE ORGANICS

All values are ug/Kg.

### Sample Identification

<u>Dilution Factor</u>	1.00	1.04	_1.43_	1.02	2.42	
Method Blank I.D.	>B4962	>B4962	>B4962	>B4962	<u>&gt;B4962</u>	
Compound	Method <u>Blank</u>		SVB 134042	SVB 136062	SVB 138082	Quantitation Limits with no <u>Dilution</u>
Chloromethane	U	U	U	U	U	10
Bromomethane	U	U	U	U	U	10
Vinyl Chloride	Ų	U	U	U	U	10
Chloroethane	U	U	U	U	U	10
Methylene Chloride	U	U	U	U	U	5
Acetone	20	16B	24B	21B	U	10
Carbon Disulfide	U	U	U	U	U	5
1,1-Dichloroethene	U	U	Ü	U	U	5 5
1,1-Dichloroethane	U	U	U	U	U	5
1,2-Dichloroethene (total)	U	U	U	U	65	5
ነloroform	U	U	U	U	U	5
<b>─</b> ,2-Dichloroethane	U	U	U	U	U	5
2-Butanone	6J	U	U	U	U	10
1,1,1-Trichloroethane	U	U	U	U	U	5
Carbon Tetrachloride	U	U	U	U	U	5
Vinyl Acetate	U	U	U	U	U	10
Bromodichloromethane	U	U	U	U	U	5
1,2-Dichloropropane	U	U	U	U	U	5
cis-1,3-Dichloropropene	· U	U	U	U	U	5
Trichloroethene	U	U	U	U	U	5
Dibromochloromethane	Ü	Ü	U	U	U	5
1,1,2-Trichloroethane	U	Ū	Ū	U	Ū	5
Benzene	Ū	Ū	Ū	Ū	Ü	5
trans-1,3-Dichloropropene	U	Ú	U	U	U	5
Bromoform	U	U	Ū	U	U	5
4-Methyl-2-pentanone	Ū	Ū	Ū	Ū	Ū	10
2-Hexanone	Ù	Ü	Ū	Ü	U	10
Tetrachloroethene	Ū	Ū	Ū	Ú	Ū	5
1,1,2,2-Tetrachloroethane	Ū	Ü	Ū	Ū	Ū	5
Toluene	Ū	Ū	Ū	Ū	Ū	5
Chlorobenzene	Ŭ	Ŭ	Ū.	Ŭ	Ü	5
Ethylbenzene	Ū	Ŭ	Ŭ	Ŭ	Ü	5
Styrene	Ŭ	Ŭ	Ū	Ŭ	Ü	5
Xylene (total)	Ü	Ü	Ŭ	Ü	Ü	5

U, J, B - See Appendix for definition.

te: Sample detection limit = quantitation limit x dilution factor.

## TABLE 2.0 30910-1527 UNISYS CORPORATION VOLATILE TENTATIVELY IDENTIFIED COMPOUNDS

Sample Identification: Method Blank >B4946

CAS#	Compound	RT	Estimated Concentration, ug/L
	None detected		
	Sample Identification:	TB 07/27/91	
CAS#	Compound	RT	Estimated Concentration, ug/L
	None detected		
	Sample Identification: Meth	od Blank >B	4962
CAS#	Compound	RT	Estimated <u>Concentration, ug/Kg</u>
	None detected		
	Sample Identification:	SVB112022	
CAS#	Compound	<u>RT</u>	Estimated <u>Concentration, ug/Kg</u>
	None detected		
	Sample Identification:	SVB114042	
CAS#	Compound	RT	Estimated <u>Concentration, ug/Kg</u>
	None detected		
, <del></del>	Sample Identification:	SVB116062	•
CAS#	Compound	RT	Estimated Concentration, ug/Kg
	None detected		

## TABLE 2.1 30910-1527 UNISYS CORPORATION VOLATILE TENTATIVELY IDENTIFIED COMPOUNDS

Sample Identification: SVB118082

CAS#	Compound None detected	<u>RT</u>	Estimated Concentration, ug/Kg		
	Sample Identification:	SVB132022			
CAS#_	Compound	RT	Estimated Concentration, uq/Kg		
91203	Naphthalene	22.58	7J		
	Sample Identification:	SVB134042			
CAS#	Compound	RT	Estimated Concentration, ug/Kg		
	None detected				
	Sample Identification:	SVB136062			
CAS#	Compound	RT	Estimated Concentration, uq/Kq		
	None detected				
	Sample Identification:	SVB138082			
CAS#	Compound	<u>RT</u>	Estimated Concentration, ug/Kg		
1634-04-4	Methyl tert butyl ether	8.66	<b>2</b> 6J		
J - See Appendix for definition.					

### APPENDIX

- U Indicates that the compound was analyzed for but not detected.
- J Indicates that the compound was analyzed for and determined to be present in the sample. The mass spectrum of the compound meets the identification criteria of the method. The concentration listed is an estimated value, which is less than the specified minimum detection limit but is greater than zero.
- B This flag is used when the analyte is found in the blanks as well as the sample. It indicates possible sample contamination and warns the data user to use caution when applying the results of this analyte.
- N Indicates that the compound was analyzed for but not requested as an analyte. Value will not be listed on tabular result sheet.
- X Matrix spike compound.
- (1) Cannot be separated from diphenylamine.
- (2) Decomposes to azobenzene. Measured and calibrated as azobenzene.
- A This flag indicates that a TIC is a suspected aldol condensation product.
- E Indicates that it exceeds calibration curve range.
- D This flag identifies all compounds identified in an analysis at a secondary dilution factor.
- C Confirmed by GC/MS.
- T Compound present in TCLP blank.



### an environmental testing company -

200 Monroe Turnpike Monroe, Connecticut 06468 (203) 261-4458 FAX (203) 268-5346

### REPORT TRANSMITTAL

REPORT NUMBER 30910-1578

DATE September 3, 1991

CLIENT

Unisys Corporation 3199 Pilot Knob MS F1B05 Eagan, MN 55121

ATTENTION Mr. Kevin Krueger

The above referenced report is enclosed. Copies of this report and supporting data will be retained in our files in the event they are required for future reference.

If there are any questions concerning this report, please do not hesitate to contact us.

Any samples submitted to our Laboratory will be retained for a maximum of sixty (60) days from receipt of this report, unless other arrangements are desired.

30910-1578
UNISYS CORPORATION
3199 Pilot Knob
MS F1B05
Eagan, Minnesota 55121

Re: Great Neck, New York

Attention: Mr. Kevin Krueger

### **PURPOSE**

Six samples and two trip blanks were submitted to IEA, Inc. by Unisys Corporation. The client requested the samples be analyzed for TCL volatile organics plus a library search for non-target compounds.

### **METHODOLOGY**

Volatile organics were determined using purge and trap GC/MS. The instrumentation used was a Tekmar Dynamic Headspace Concentrator interfaced with a Hewlett-Packard Model 5995C GC/MS/DS.

### DISCUSSION

<u>Volatile Organics</u> - The laboratory followed USEPA CLP-SOW Document #OLMO1.0 for the GC/MS calibration criteria.

Samples SVB-10 20-22 and SVB-10 40-42 were analyzed twice within NYSDEC holding times but the data was unacceptable due to retention time shifts. The samples were reanalyzed on 08/13/91 (one day past CLP holding times) but the method blank had acetone at 150 ppb (out of criteria). Acetone was also found in the samples at 100 ppb and 190 ppb respectively. The samples were reanalyzed again on 08/15/91 with no problems. The client was contacted and requested the runs from 08/13/91 and 08/15/91 be reported. The samples run on 08/15/91 have been flagged with the suffix "RE".

Sample TB 07/30/91 was analyzed out of NYSDEC holding time and 3 days past CLP holding time on 08/15/91. This run has been reported. Both trip blanks were analyzed by CLP 2/88 criteria, (NYSDEC '89 criteria could not be achieved).

Sample SVB-15 40-42 was analyzed 2 hours past NYSDEC holding time but within CLP holding time. The client requested the run be reported.

TIC's detected in method blank (VBLKAA) were caused by system carryover. These TIC's were not detected in any associated sample.

## **RESULTS**

The results are presented in the following Tables. Also enclosed is the data package containing all relevant data.

Prepared by

effrey C. Curran abonatory Manager

JCC/mt

The liability of IEA, Inc. is limited to the actual dollar value of this project.

## TABLE 1.0 30910-1578 UNISYS CORPORATION **EPA TCL VOLATILE ORGANICS**

All values are ug/L.

<u>Dilution Factor</u>	1.0	_1.0	
Method Blank I.D.	<u>&gt;G5863</u>	<u>&gt;G5863</u>	
	Method	ТВ	Quantitation Limits with no
<u>Compound</u>	<u>Blank</u>	08/07/91	<u>Dilution</u>
Chloromethane	Ų	Ų	10
Bromomethane	Ü	U	10
Vinyl Chloride	U	U	10
Chloroethane	Ų	U	10
Methylene Chloride	U	Ų	5
Acetone	U	Ų	10
Carbon Disulfide	Ų	Ų .	5
1,1-Dichloroethene	Ų	U	5
1,1-Dichloroethane	U	Ų	5
1,2-Dichloroethene (total) Chloroform	U U	Ų	5 5
1,2-Dichloroethane	U	U U	5 5
2-Butanone	Ü	Ü	10
1,1,1-Trichloroethane	Ü	Ŭ	5
Carbon Tetrachloride	Ü	Ü	5
Vinyl Acetate	Ŭ	Ŭ	10
Bromodichloromethane	Ü	Ŭ	5
1,2-Dichloropropane	Ŭ	Ŭ	5
cis-1,3-Dichloropropene	Ŭ	ŭ	5
Trichloroethene	Ü	Ŭ	5
Dibromochloromethane	Ŭ	Ŭ	5
1,1,2-Trichloroethane	Ü	Ŭ	5
Benzene	Ü	Ü	5
trans-1,3-Dichloropropene	Ü	Ü	5 5
Bromoform	U	U	5
4-Methyl-2-pentanone	U	U	10
2-Hexanone	U	U	10
Tetrachloroethene	U	U	5
1,1,2,2-Tetrachloroethane	U	U	5
Toluene	U	- U -	5
Chlorobenzene	Ų	U	5
Ethylbenzene	Ü	U	5 5
Styrene (tatal)	Ų	U	5
Xylene (total)	U	U	5

U - See Appendix for definition. Note: Sample detection limit = quantitation limit x dilution factor.

## TABLE 1.1 30910-1578 UNISYS CORPORATION **EPA TCL VOLATILE ORGANICS**

All values are ug/L.

Sample Identification

<u>Dilution Factor</u>	1.0	1.0	
Method Blank I.D.	<u>&gt;G5881</u>	<u>&gt;G5881</u>	
Compound	Method <u>Blank</u>	TB <u>07/30/91</u>	Quantitation Limits with no <u>Dilu</u> tion
Chloromethane	U	U	10
Bromomethane	U	U	10
Vinyl Chloride	U	U	10
Chloroethane	U	U	10
Methylene Chloride	U	U	5
Acetone	U	U	10
Carbon Disulfide	U	U	5
1,1-Dichloroethene	U	U	5
1,1-Dichloroethane	U	U	5
1,2-Dichloroethene (total)	U	U	5 5
Chloroform	Ū	Ū	5
1,2-Dichloroethane	U	U	5
2-Butanone	U	U	10
1,1,1-Trichloroethane	U	U	5
Carbon Tetrachloride	U	U	5
Vinyl Acetate	U	U	10
Bromodichloromethane	U	U	5
1,2-Dichloropropane	Ū	Ü	5
cis-1,3-Dichloropropene	Ū	Ü	5
Trichloroethene	Ŭ	Ü	5
Dibromochloromethane	Ŭ	Ŭ	5
1,1,2-Trichloroethane	Ŭ	Ü	5
Benzene	Ŭ	Ŭ	5
trans-1,3-Dichloropropene	Ü	Ü	5
Bromoform	Ŭ	Ŭ	5
4-Methyl-2-pentanone	Ŭ	Ŭ	10
2-Hexanone	Ŭ	Ū	10
Tetrachloroethene	Ū	Ŭ	5
1,1,2,2-Tetrachloroethane	Ü	Ü	5
Toluene	Ŭ	· Ŭ	5
Chlorobenzene	Ŭ	Ŭ	5
Ethylbenzene	Ŭ	Ŭ	5
Styrene	Ŭ	Ŭ	5
Xylene (total)	Ŭ	Ü	5
- · · · · · · · · · · · · · · · · · · ·	_	-	•

U ~ See Appendix for definition.
Note: Sample detection limit = quantitation limit x dilution factor.

## TABLE 1.2 30910-1578 UNISYS CORPORATION **EPA TCL VOLATILE ORGANICS**

All values are ug/Kg.

<u>Dilution Factor</u>	1.00	1.04	1.05		
Method Blank I.D.	<u>&gt;A7247</u>	<u>&gt;A7247</u>	<u>&gt;A</u> 7247		
	Method	SVB-10	SVB-10		Quantitation Limits with no
<u>Compound</u>	<u>Blank</u>	20-22			<u>Dilution</u>
Chloromethane	U	U	U		10
Bromomethane	U	U	Ų		10
Vinyl Chloride Chloroethane	U U	U U	U U		10
Methylene Chloride	U	ა ვე	ა ვე		10 5
Acetone	150	100B	190B		10
Carbon Disulfide	U	U	U		5
1,1-Dichloroethene	Ŭ	Ü	Ŭ	•	5
1,1-Dichloroethane	Ü	Ŭ	Ŭ		5
1,2-Dichloroethene (total)	Ŭ	2J	10		5 5 5 5
lloroform	Ŭ	บ	Ū		5
1,2-Dichloroethane	Ŭ	Ŭ	Ū		5
2-Butanone	0.8J	Ū	<b>2</b> JB		10
1,1,1-Trichloroethane	U	V	U		5
Carbon Tetrachloride	U	U	U		5
Vinyl Acetate	U	U	U		10
Bromodichloromethane	U	U	U		5
1,2-Dichloropropane	U	U	U		5 5 5 5
cis-1,3-Dichloropropene	U	U	U		5
Trichloroethene	U	U	U		5
Dibromochloromethane	U	U	U		5_
1,1,2-Trichloroethane	U	U	Ü		5 5 5
Benzene	Ų	Ü	U		5
trans-1,3-Dichloropropene Bromoform	U U	U U	U U		5 5
4-Methyl-2-pentanone	U	U	U		10
2-Hexanone	Ü	Ü	Ü		10
Tetrachloroethene	Ü	2J	5		5
1,1,2,2-Tetrachloroethane	Ŭ	U	U		5
Toluene	Ū	Ŭ	Ū	•	5
Chlorobenzene	Ŭ	Ŭ	Ŭ		5
Ethylbenzene	U	Ü	Ū		
Styrene	U	U	U		5 5
Xylene (total)	0.8J	U	U		5

<sup>&#</sup>x27;, J, B - See Appendix for definition. total one one one of the property o

## TABLE 1.3 30910-1578 **UNISYS CORPORATION EPA TCL VOLATILE ORGANICS**

All values are ug/Kg.

<u>Dilution Factor</u>	1.00	1.03	1.02	1.19	1.32	
Method Blank I.D.	<u>&gt;A7272</u>	<u>&gt;A7272</u>	<u>&gt;A7272</u>	<u>&gt;A7272</u>	<u>&gt;A7272</u>	Quantitation
<u>Compound</u>				SVB-15 80-82	SVB-15 40-42	Limits with noDilution
Chloromethane	U	U	U	U	U	10
Bromomethane	U	U	U	U	U	10
Vinyl Chloride	U	U	U	U	U	10
Chloroethane	U	U	U	U	U	10
Methylene Chloride	3J	1JB	1JB	U	U	5
Acetone	15	7JB	7JB	U	17B	10
Carbon Disulfide	U	U	U	U	. <b>U</b>	5
1,1-Dichloroethene	U	U	U	U	U	5
1,1-Dichloroethane	U	IJ	U	U	U	5
1,2-Dichloroethene (total)	U	U	U	U	U	5
`hloroform	U	U	U	U	U	5
■1,2-Dichloroethane	U	U	U	U	U	5
2-Butanone	2J	Ü	U	U	U	10
1,1,1-Trichloroethane	U	U	U	U	U	5
Carbon Tetrachloride	U	U	IJ	U	U	5
Vinyl Acetate	U	U	U	U	U	10
Bromodichloromethane	U	U	U	U	U	5
1,2-Dichloropropane	U	U	U	U	U	5
cis-1,3-Dichloropropene	U	U	IJ	U	U	5
Trichloroethene	U	U	U	U	U	5
Dibromochloromethane	U	U	U	U	U	5
1,1,2-Trichloroethane	U	U	U	U	U	5
Benzene	U	U	U	U	U	5 5
trans-1,3-Dichloropropene	U	U	U	U	U	
Bromoform	U	U	U	U	U	5
4-Methyl-2-pentanone	0.6J	U	U	U	U	10
2-Hexanone	U	U	U	U	U	10
Tetrachloroethene	U	U	U	U	U	5
1,1,2,2-Tetrachloroethane	U	U	IJ	U	U	5
Toluene	U	U	U	U	U_	5
Chlorobenzene	U	U	U	U	U	5
Ethylbenzene	U	U	U	U	U	5
Styrene	U	U	U	U	U	5 5 5
Xylene (total)	U	U	U	U	U	5

U, J, B - See Appendix for definition. \_lote: Sample detection limit = quantitation limit x dilution factor.

## TABLE 1.4 30910-1578 UNISYS CORPORATION **EPA TCL VOLATILE ORGANICS**

All values are ug/Kg.

<u>Dilution Factor</u>	1.00	1.03	1.04	
Method Blank I.D.	<u>&gt;B5116</u>	>B5116 SVB-10	>B5116 SVB-10	Quantitation
Compound	Method <u>Blank</u>	20-22 RE	40-42 RE	Limits with no <u>Dilution</u>
Chloromethane	U	U	U	10
Bromomethane	U	U	U	. 10
Vinyl Chloride	U	U	U	10
Chloroethane	U	U	Ú	10
Methylene Chloride	<b>4</b> J	U	U	5
Acetone	<b>9</b> J	13B	5JB	10
Carbon Disulfide	U	U	U	. 5
1,1-Dichloroethene	U	U	U	5
1,1-Dichloroethane	U	U	U	
1,2-Dichloroethene (total)	Ü	U	U	5 5 5
^hloroform	Ū	Ü	U	5
,2-Dichloroethane	Ú	Ü	U	5
2-Butanone	0.7J	U	U	10
1,1,1-Trichloroethane	U	U	U	5
Carbon Tetrachloride	U	U	U	5
Vinyl Acetate	U	U	U	10
Bromodichloromethane	U	U	U	5
1,2-Dichloropropane	U	U	Ú	5
cis-1,3-Dichloropropene	U	U	U	5
Trichĺoroethene .	U	U	U	5 5 5 5
Dibromochloromethane	U	Ū	U	5
1,1,2-Trichloroethane	U	Ū	Ü	5
Benzene	U	U	U	5
trans-1,3-Dichloropropene	U	U	U	5
Bromoform	U	U	U	5
4-Methyl-2-pentanone	U	U	U	10
2-Hexanone	U	U	U	10
Tetrachloroethene	U	U	U	5
1,1,2,2-Tetrachloroethane	U	U	U	5 5
Toluene	U	U	U	<u>_</u> 5
Chlorobenzene	U	U	U	5
Ethylbenzene	U	U	U	5
Styrene	U	U	U	5
Xylene (total)	2J	U	U	5

U, J, B - See Appendix for definition. ote: Sample detection limit  $\Rightarrow$  quantitation limit x dilution factor.

# TABLE 2.0 30910-1578 UNISYS CORPORATION VOLATILE TENTATIVELY IDENTIFIED COMPOUNDS

Sample Identification: Method Blank >G5863

	Sample Identification: Method	I Blank >G	5863
CAS#	Compound	RT	Estimated Concentration, ug/L
	None detected		
	Sample Identification: TB	08/07/91	
CAS#	Compound	RT	Estimated Concentration, ug/L
87683	1,1,2,3,4,4-Hexachloro-1,3- butadiene	27.15	<b>59</b> J
91203	Unknown trichlorobenzene isomer Unknown trichlorobenzene isomer Naphthalene	27.68 26.95 27.30	38J 20J 16J
	Sample Identification: Method	Blank >G	5881
CAS#	Compound	RT	Estimated Concentration, ug/L
	None detected		
	Sample Identification: TB	07/30/91	
CAS#	Compound	RT	Estimated Concentration, ug/L
	None detected		
	Sample Identification: Method	Blank >A	7247
CAS#	Compound	RT	Estimated <u>Concentration, ug/Kg</u>
96140	3-Methyl-pentane	9.29	11J
J - See A	ppendix for definition.		

# TABLE 2.1 30910-1578 UNISYS CORPORATION VOLATILE TENTATIVELY IDENTIFIED COMPOUNDS

Sample Identification: SVB-10 20-22

CAS#	Compound	<u>RT</u>	Estimated Concentration, ug/Kg
	Unknown branched alkane	24.09	10J
	Unknown C, alkyl benzene	25.75	10J
	Unknown cycloalkane	26.34	10J
	Unknown cycloalkane	24.55	9J
	Unknown cycloalkane	22.37	8J
	Unknown cycloalkane	23.09	8J
	Unknown alkane	25.95 24.42	7J
	Unknown branched alkane	24.42 23.54	7J 6J
	Unknown branched cycloalkane Unknown branched alkane	22.14	6J
	onknown branched arkane	22.14	00
	Sample Identification: S	VB-10 40-42	
			Estimated
CAS#	Compound	<u>RT</u>	Concentration, ug/Kg
	None detected		
	Sample Identification: Metho	od Blank >A7	272
			Estimated
CAS#	<u>Compound</u>	RT	Concentration, ug/Kg
	Unknown	21.65	62J
	Unknown isomer of trichlorobenzene	19.18	58J
91203	Naphthalene	21.19	<b>52</b> J
	Comming Vilonities and the continue of	WD 15 00 00	
	Sample Identification: S	AR-12 50-55	
	•		Estimated
CAS#	Compound	RT	Concentration, ug/Kg
	None detected		•

 $\ensuremath{\mathsf{J}}$  - See Appendix for definition.

## TABLE 2.2 30910-1578 UNISYS CORPORATION VOLATILE TENTATIVELY IDENTIFIED COMPOUNDS

Sample Identification: SVB-15 60-62

	Sample Identification: S	SVB-15 60-62	,
CAS#	Compound	. <u>RT</u>	Estimated Concentration, ug/Kg
	None detected		
	Sample Identification: S	SVB-15 80-82	
CAC#	Compound	RT	Estimated
CAS#	Compound	_ <u>KI</u>	Concentration, uq/Kq
	None detected		
	Sample Identification: S	SVB-15 40-42	
			Estimated
CAS#	Compound	<u>RT</u>	Concentration, ug/Kg
	None detected		
	Sample Identification: Meth	od Blank >B5	5116
			Estimated
<u>CAS#</u>	<u>Compound</u>	<u>RT</u>	Concentration, ug/Kg
	None detected		
	Sample Identification: SV	B-10 20-22 F	RE
			Estimated
CAS#	Compound	<u>RT</u>	Concentration, uq/Kq
	Unknown C, alkyl benzene	26.51	23J
	Unknown branched alkene	25.86	16J
2958761			
	Decahydro-2-methyl-naphthalene	26.06	14J
	Unknown alkyl benzene	25.24	<b>-</b> 12J
	Unknown alkyl benzene Unknown alkane	25.24 25.44	• 12J 8J
	Unknown alkyl benzene Unknown alkane Unknown alkane	25.24 25.44 23.52	* 12J 8J 8J
	Unknown alkyl benzene Unknown alkane Unknown alkane Unknown alkylbenzene	25.24 25.44 23.52 24.95	* 12J 8J 8J <b>7</b> J
	Unknown alkyl benzene Unknown alkane Unknown alkane Unknown alkylbenzene Unknown decahydronaphthalene isome	25.24 25.44 23.52 24.95 r 24.79	- 12J 8J 8J 7J 7J
	Unknown alkyl benzene Unknown alkane Unknown alkane Unknown alkylbenzene	25.24 25.44 23.52 24.95	* 12J 8J 8J <b>7</b> J

 $<sup>{\</sup>sf J}$  - See Appendix for definition.

# TABLE 2.3 30910-1578 UNISYS CORPORATION VOLATILE TENTATIVELY IDENTIFIED COMPOUNDS

Sample Identification: SVB-10 40-42RE

		•	Estimated
CAS#	Compound	RT	Concentration, ug/Kg
	i		
	None detected		

#### APPENDIX

- U Indicates that the compound was analyzed for but not detected.
- J Indicates that the compound was analyzed for and determined to be present in the sample. The mass spectrum of the compound meets the identification criteria of the method. The concentration listed is an estimated value, which is less than the specified minimum detection limit but is greater than zero.
- B This flag is used when the analyte is found in the blanks as well as the sample. It indicates possible sample contamination and warns the data user to use caution when applying the results of this analyte.
- N Indicates that the compound was analyzed for but not requested as an analyte. Value will not be listed on tabular result sheet.
- X Matrix spike compound.
- (1) Cannot be separated from diphenylamine.
- (2) Decomposes to azobenzene. Measured and calibrated as azobenzene.
- A This flag indicates that a TIC is a suspected aldol condensation product.
- E Indicates that it exceeds calibration curve range.
- D This flag identifies all compounds identified in an analysis at a secondary dilution factor.
- C Confirmed by GC/MS.
- T Compound present in TCLP blank.



an environmental testing company -

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#### REPORT TRANSMITTAL

REPORT NUMBER 30910-1662

DATE September 17, 1991\_

CLIENT

Unisys Corporation 3199 Pilot Knob MS F1B05 Eagan, MN 55121

ATTENTION Mr. Kevin Krueger

The above referenced report is enclosed. Copies of this report and supporting data will be retained in our files in the event they are required for future reference.

If there are any questions concerning this report, please do not hesitate to contact us.

Any samples submitted to our Laboratory will be retained for a maximum of thirty (30) days from receipt of this report, unless other arrangements are desired.

30910-1662 UNISYS CORPORATION 3199 Pilot Knob MS F1B05 Eagan, Minnesota 55121

Re: Great Neck, New York

Attention: Mr. Kevin Krueger

#### **PURPOSE**

Eight samples and two trip blanks collected on August 14-16, 1991 were submitted to IEA, Inc. by Unisys Corporation. The client requested the samples be analyzed for TCL volatile organics plus a library search for non-target compounds.

#### <u>METHODOLOGY</u>

Volatile organics were determined using purge and trap GC/MS. The instrumentation used was a Tekmar Dynamic Headspace Concentrator interfaced with a Hewlett-Packard Model 5995C GC/MS/DS.

#### **DISCUSSION**

<u>Volatile Organics</u> - The laboratory followed USEPA CLP-SOW Document #OLMO1.6 for the GC/MS calibration criteria.

The 08/15/91 trip blank was analyzed on 08/23/91 (one day past NYSDEC ASP holding time) due to instrumentation problems.

The low level water initial calibration on Instrument G on 08/22/91 contained five percent RSD's above the 20.5 percent criteria. This exceeds the 0LM01.6 criteria by three compounds; however, the trip blanks were analyzed under this curve due to holding time considerations. The continuing calibration (>G6026) had no compounds out of criteria when compared to the above curve.

### **RESULTS**

The results are presented in the following Tables. Also enclosed is the data package containing all relevant data.

Prepared by:

Curran Manager

JCC/mt

The liability of IEA, Inc. is limited to the actual dollar value of this project.

## TABLE 1.0 30910-1662 UNISYS CORPORATION **EPA TCL VOLATILE ORGANICS**

All values are ug/L.

Sample Identification

<u>Dilution Factor</u>	1.0_	1.0	1.0	
Method Blank I.D.	<u>&gt;G6027</u>	<u>&gt;G6027</u>	<u>&gt;G6027</u>	0
<u>Compound</u>	Method <u>Blank</u>	TB <u>08/15/91</u>	TB <u>08/16/91</u>	Quantitation Limits with no <u>Dilution</u>
Chloromethane	U	U	U	10
Bromomethane	U	U	U	10
Vinyl Chloride	U	U	U	10
Chloroethane	U	U	U	10
Methylene Chloride	<b>2</b> J	U	U	5
Acetone	12	U	12B	10
Carbon Disulfide	U	U	U	5
1,1-Dichloroethene	U	U	U ·	5
1,1-Dichloroethane	Ü	U	U	5
1,2-Dichloroethene (total)	Ū	Ū	Ü	5
Chloroform	Ŭ	Ŭ	Ü	5
1,2-Dichloroethane	Ŭ	Ü	Ü	5
2-Butanone	Ŭ	Ŭ	Ü	10
1,1,1-Trichloroethane	Ŭ	Ŭ	Ü	5
Carbon Tetrachloride	ŭ	Ŭ	Ü	5
Vinyl Acetate	ŭ	Ŭ	Ü	10
Bromodichloromethane	ŭ	Ŭ	Ü	5
1,2-Dichloropropane	Ŭ	Ŭ	Ü	5
cis-1,3-Dichloropropene	Ŭ	Ů	Ŭ	5
Trichloroethene	Ŭ	Ŭ	Ü	5
Dibromochloromethane	Ü	Ŭ	Ŭ٠、	5
1,1,2-Trichloroethane	Ŭ	Ŭ	Ŭ	5
Benzene	Ŭ	ŭ	ŭ	5
trans-1,3-Dichloropropene	Ŭ	บ	Ü	5
Bromoform	Ŭ	Ŭ	Ü	5
4-Methyl-2-pentanone	Ŭ	Ŭ	Ü	10
2-Hexanone	Ŭ	Ŭ	Ŭ	10
Tetrachloroethene	Ŭ	Ŭ	Ŭ	5
1,1,2,2-Tetrachloroethane	Ü	Ŭ	Ü	5
Toluene	0.7J	Ü	Ŭ	
Chlorobenzene	Ü	Ŭ	Ŭ =	5 5
Ethylbenzene	Ü	Ü	Ü	5
Styrene	U	Ü	Ü	5
Xylene (total)	Ü	Ü	Ü	5
Afrene (Lucai)	U	U	U	<b>5</b>

U, J, B - See Appendix for definition. Note: Sample detection limit = quantitation limit  ${\bf x}$  dilution factor.

## TABLE 1.1 30910-1662 UNISYS CORPORATION EPA TCL VOLATILE ORGANICS

All values are ug/Kg.

## Sample Identification

<u>Dilution Factor</u>	_1.00	1.20	1.04	1.11	1.05	1.05	
Method Blank I.D.	<u>&gt;B5234</u>	<u>&gt;B5234</u>	<u>&gt;B5234</u>	<u>&gt;B5234</u>	<u>&gt;B5234</u>	<u>&gt;B5234</u>	0
Compound			SVB-16 40-42'				Quantitation Limits with no <u>Dilution</u>
Chloromethane	U	U	U	U	U	U	10
Bromomethane	U	Ü	U	U	U	U	10
Vinyl Chloride	U	U	U	U	U	U	10
Chloroethane	U	U	U	U	U	U	10
Methylene Chloride	Ų	<b>5</b> J	6	6	7	4J	5
Acetone	<b>5</b> J	<b>9</b> JB	13B	13B	19B	26B	10
Carbon Disulfide	U	U	U	U	U	U	5
1,1-Dichloroethene	U	U	U	U	Ú .	U	5 5
1,1-Dichloroethane	U	U	U	U	U	U	5
1,2-Dichloroethene (total)	U	30	U	U	U	U	5
Chloroform	U	U	U	U	U	U	5
l,2-Dichloroethane	U	U	U	U	U	U	5
2-Butanone	U	U	U	U	U	U	10
1,1,1-Trichloroethane	U	U	U	U	U	U	5
Carbon Tetrachloride	U	U	U	U	U	Ü	5
Vinyl Acetate	U	U	U	U	U	U	10
Bromodichloromethane	U	U	U	U	U	U	5
1,2-Dichloropropane	U	U	U	U	U	U	5
cis-1,3-Dichloropropene	U	IJ	U	U	U	U	5
Trichloroethene	U	<b>2</b> J	U	U	U	U	5
Dibromochloromethane	U	U	IJ	U	A.	U	5
1,1,2-Trichloroethane	υ	U	U	U	U	Ų	5
Benzene	U	U	U	U	U	U	5
trans-1,3-Dichloropropene	U	U	U	U	U	U	5
Bromoform	U	U	U	U	U	U	5
4-Methyl-2-pentanone	U	U	U	U	U	U	10
2-Hexanone	U	U	U	U	U	U	10
Tetrachloroethene	U	2J	U	U	U	U	5
1,1,2,2-Tetrachloroethane	U	U	U	U	U	U	5
Toluene	U	U -	U	U	U	U	5
Chlorobenzene	U	U	U .	U	_	• U	5
Ethylbenzene	Ü	U	U	U	U	IJ	5
Styrene	U	U	U	ป	U	U	5
Xylene (total)	U	U	U	17	U	U	5

U, J, B - See Appendix for definition. Note: Sample detection limit = quantitation limit  $\mathbf x$  dilution factor.

## TABLE 1.2 30910-1662 UNISYS CORPORATION **EPA TCL VOLATILE ORGANICS**

All values are ug/Kg.

## Sample Identification

<u>Dilution Factor</u>	1.00	1.14	1.23	1.04	1.20	1.20	
Method Blank I.D.	<u>&gt;B5234</u>	<u>&gt;B5234</u>	<u>&gt;B5234</u>	<u>&gt;B5234</u>	>B5234 SVB-16	>B5234 SVB-16	Quantitation
<u>Compound</u>			SVB-14 80-82'		80-82'	80-82'	Limits with no <u>Dilution</u>
Chloromethane	U	U	U	U	U	U	10
Bromomethane	U	U	U	U	U	U	10
Vinyl Chloride	U	U	U	U	U	U	10
Chloroethane	U	U	U	U	U	U	10
Methylene Chloride	U	<b>4</b> J	U	<b>4</b> J	9	3ე	5
Acetone	5J	40B	. U	29B	22B	7JB	10
Carbon Disulfide	U	U	U	U	U	U	5
1,1-Dichloroethene	U	U	U	U	70X	70X	5
1,1-Dichloroethane	U	U	U	U	U	U	5
1,2-Dichloroethene (total)	U	U	U	U	19	8	5
Chloroform	U	U	U	U	U	U	5
_1,2-Dichloroethane	U	U	U	U	U	U	5
2-Butanone	U	U	U	U	U	U	10
1,1,1-Trichloroethane	U	U	U	U	U	U	5
Carbon Tetrachloride	U	U	U	U	U	U	5
Vinyl Acetate	U	U	U	U	U	U	10
Bromodichloromethane	U	U	U	U	U	U	5
1,2-Dichloropropane	U	U	U	U	U	U	5
cis-1,3-Dichloropropene	U	U	U	U	U	U	5
Trichĺoroethene	U	U	U	U	61X	60X	5
Dibromochloromethane	IJ	U	U	U	Ų.	U	5
1,1,2-Trichloroethane	U	U	U	U	U	U	5
Benzene	U	U	U	U	66X	66X	5
trans-1,3-Dichloropropene	U	U	U	U	U	U	5
Bromoform	U	U	U	U	U	U	5
4-Methyl-2-pentanone	U	U	U	U	U	U	10
2-Hexanone	U	U	U	U	U	U	10
Tetrachloroethene	U	U	U	U	U	U	5
1,1,2,2-Tetrachloroethane	IJ	U	U	U	U	U	5
Toluene	U	U	U	U	64X	64X	5 5
Chlorobenzene	U	U	U -	U	60X 🛥		5
Ethylbenzene	U	U	U	U	U	U	5
Styrene	U	U	U	U	U	U	5
Xylene (total)	U	IJ	15	U	U	U	5
·							

U, J, B, X - See Appendix for definition. Note: Sample detection limit = quantitation limit  ${\bf x}$  dilution factor.

## TABLE 2.0 30910-1662 UNISYS CORPORATION VOLATILE TENTATIVELY IDENTIFIED COMPOUNDS

Sample Identification: Method Blank >G6027

CAS#	Compound None detected	<u>RT</u>	Estimated Concentration, ug/L
	Sample Identification: TB	08/15/91	
CAS#	<u>Compound</u>	RT	Estimated <u>Concentration, uq/L</u>
	None detected		
	none detected		
	Sample Identification: TB	08/16/91	Fakimakad
CAS#	Compound	RT	Estimated Concentration, ug/L
CA3#	Compound		concentration, ug/L
	None detected		
	Sample Identification: Method	Blank >B	5234
			Estimated
<u>CAS#</u>	Compound	<u>RT</u>	Concentration, ug/Kg
	Unknown alkane Unknown alkane Unknown	26.29 25.80 22.80	16J 10J 6J
	Sample Identification: SVB-	-16 80-82′	
CAS#	<u>Compound</u>	RT	Estimated Concentration, ug/Kg
	Unknown alkane	25.17	10J
	Unknown alkane	23.47	<b>-</b> 8J
	Unknown alkane	25.75	8JB
	Unknown alkane	24.06	<b>7</b> J
	Unknown alkane	26.24	7JB

J, B - See Appendix for definition.

## TABLE 2.1 30910-1662 UNISYS CORPORATION VOLATILE TENTATIVELY IDENTIFIED COMPOUNDS

Sample Identification: SVB-16 40-42'

CAS#	Compound Unknown alkane Unknown Unknown alkane Unknown	RT	Estimated Concentration, ug/Kg  8JB 6J 6J 6J 6J
	Sample Identification:	SVB-16 60-62'	
CAS#	Compound	RT	Estimated Concentration, ug/Kg
	Unknown alkene Unknown alkene	26.30 23.51	14J 12J
	Unknown	25.78	8J
	Unknown alkane	24.19	8J
	Unknown	25.26	6J
	Sample Identification:	SVB-16 20-22'	•
CAS#	Compound	<u>RT</u>	Estimated Concentration, ug/Kg
	Unknown alkane	26.50	13JB
	Unknown alkane	26.07	8J
	Unknown alkene	24.22	8J 7J
	Unknown ketone Unknown	23.01 23.66	70 5J
	Sample Identification:		30
CAS#	Compound	RT	Estimated <u>Concentration, ug/Kg</u>
	Unknown alkane	23.50	<b>-</b> 7J

J, B - See Appendix for definition.

## TABLE 2.2 30910-1662 UNISYS CORPORATION VOLATILE TENTATIVELY IDENTIFIED COMPOUNDS

Sample Identification: SVB-14 40-42'

CAS#	Compound	RT	Estimated Concentration, ug/Kg
	Unknown alkane Unknown C <sub>3</sub> alkyl benzene	24.93 23.50	26J 22J
	Unknown alkane Unknown alkane	23.01 20.54	22J 16J
	Unknown branched alkane	22.23	15J
	Unknown alkene Unknown C <sub>r</sub> alkyl benzene	21.74 26.14	12J 12J
	Unknown C <sub>5</sub> alkyl benzene Unknown C <sub>4</sub> alkyl benzene Unknown cycloalkane	25.22 24.02	11J 11J
	Unknown	24.18	10J
	Sample Identification	n: SVB-14 80-82	,
CAS#	Compound	<u></u>	Estimated Concentration, ug/Kg
	Unknown C <sub>3</sub> alkyl benzene Unknown	23.51 24.65	7J 7J
	Sample Identification	n: SVB-14 20-22	,
<u>CA</u> S#	Compound		Estimated Concentration, ug/Kg
	Unknown C <sub>3</sub> alkyl benzene Unknown	23.50 26.27	7J 6J

 $<sup>\</sup>ensuremath{\mathbf{J}}$  - See Appendix for definition.

#### APPENDIX

- U Indicates that the compound was analyzed for but not detected.
- J Indicates that the compound was analyzed for and determined to be present in the sample. The mass spectrum of the compound meets the identification criteria of the method. The concentration listed is an estimated value, which is less than the specified minimum detection limit but is greater than zero.
- B This flag is used when the analyte is found in the blanks as well as the sample. It indicates possible sample contamination and warns the data user to use caution when applying the results of this analyte.
- N Indicates that the compound was analyzed for but not requested as an analyte. Value will not be listed on tabular result sheet.
- X Matrix spike compound.
- (1) Cannot be separated.
- (2) Decomposes to azobenzene. Measured and calibrated as azobenzene.
- A This flag indicates that a TIC is a suspected aldol condensation product.
- E Indicates that it exceeds calibration curve range.
- D This flag identifies all compounds identified in an analysis at a secondary dilution factor.
- C Confirmed by GC/MS.
- T Compound present in TCLP blank.



#### an environmental testing company -

200 Monroe Turnpike Monroe, Connecticut 06468 (203) 261-4458 FAX (203) 268-5346

## REPORT TRANSMITTAL

REPORT NUMBER 30910-1721

DATE\_\_\_\_\_October 2, 1991

CLIENT

Unisys Corporation 3199 Pilot Knob MS F1B05

Eagan, Minnesota 55121

ATTENTION <u>Mr. Kevin Krueger</u>

The above referenced report is enclosed. Copies of this report and supporting data will be retained in our files in the event they are required for future reference.

If there are any questions concerning this report, please do not hesitate to contact us.

Any samples submitted to our Laboratory will be retained for a maximum of thirty (30) days from receipt of this report, unless other arrangements are desired.

30910-1721 UNISYS CORPORATION 3199 Pilot Knob MS F1B05 Eagan, Minnesota 55121

Re: Great Neck, New York

Attention: Mr. Kevin Krueger

#### **PURPOSE**

Thirteen samples collected on August 21, 1991 were submitted to IEA, Inc. by Unisys Corporation. The client requested the samples be analyzed for the parameters listed in Table 1.0.

#### METHODOLOGY

Volatile organics were determined using purge and trap GC/MS. The instrumentation used was a Tekmar Dynamic Headspace Concentrator interfaced with a Hewlett-Packard Model 5995C GC/MS/DS.

Semi-volatile organics were determined using capillary GC/MS. The instrumentation used was a Hewlett-Packard Model 5890 gas chromatograph interfaced with a Model 5970 Mass Selective Detector.

Pesticides and polychlorinated biphenyls (PCB's) were determined using GC/ECD. The instrumentation used was a HP Model 5890 gas chromatograph equipped with an electron capture detector ( $Ni^{63}$ ).

Metals were determined by ICP using either a JA61 simultaneous ICAP or a PE6500-XR sequential ICP. Graphite furnace elements were determined using either a PEZ5100 or a PEZ3030 GFAAS. Mercury was determined by the cold vapor technique utilizing the Spectro Products Model HG-4 mercury analyzer.

Cyanide was determined colorimetrically after preliminary distillation.

All analyses were conducted according to NYSDEC '89 Protocols.

#### **DISCUSSION**

<u>Miscellaneous</u> - Twelve samples were received at IEA, Inc. on 08/23/91, but due to an indistinct chain-of-custody the laboratory logged in thirteen samples.

Numerous attempts were made to get a revised chain-of-custody, however a response was not made therefore the laboratory took the initiative to log in the samples based on the sample bottle labels. Two bottle sets were received for sample SVB-17 40-42' for volatiles which was written on the original chain-of-custody. The laboratory interpreted it to be two separate samples for two reasons: 1) the sample ID's being written on separate lines on the original chain-of-custody; and 2) the sample ID's were written under two sample bottle sets.

The laboratory received the revised chain-of-custody on 09/13/91 with sample SVB-17 40-42' listed as only one sample. The analyses had already been performed (holding times were due on 08/30/91) therefore two sample results were reported for SVB-17 40-42. The results were reported under the following sample ID's: SVB-17 40-42' 002 and SVB-17 40-42' 011.

<u>Volatile Organics</u> - The initial calibration on instrument G on 08/22/91 for low level water did not meet OLM01.3 SOW calibration criteria. Four compounds exceeded the 20.5 percent RSD criteria (only two are permitted to be outside criteria).

Dibromochloromethane - 27.8% RSD 1,1,2,2-Tetrachloroethane - 21.2% RSD Xylene (total) - 27.1% RSD Bromofluorobenzene - 26.9% RSD

The target compounds listed above were not detected in any of the samples analyzed under this initial calibration.

Samples SVB-17 60-62', SVB-17 40-42' and SVB-17B 20-22' were analyzed within holding time as low level soils, however results indicated reanalysis at high medium level soil dilutions was necessary. A medium level analysis was performed for these samples within holding time but were not acceptable at their dilutions. Samples SVB-17 60-62' and SVB-17B 20-22' were successfully analyzed on 08/31/91 (eight days from sample receipt) and SVB-17 40-42' was run on 09/03/91 (eleven days from sample receipt). The client requested both sets of medium level soil data only be reported.

Sample SVB-17 80-82' was analyzed within holding time as a medium level soil, however internal standard areas were below 50 percent of the calibration standard and surrogate recoveries were diluted out. A reanalysis was performed one day past NYSDEC '89 holding time on 08/31/91 and results proved matrix interference of internal standard area recoveries. Both sets of data were reported.

Sample SVB-17 40-42' was originally analyzed within holding time at a medium level soil dilution for acetone, however the analysis proved to be over diluted. A reanalysis took place eleven days after sample receipt at the proper dilution for acetone. The QC was also run out of holding time for this sample. Both sets of data for SVB-17 40-42' were reported.

The laboratory followed the USEPA CLP-SOW document #OLMO1.0 for the GC/MS calibration criteria.

<u>Semi-Volatile Organics</u> - Sample SVB-17 12-20' exhibited suppression of the internal standard phenanthrene- $d_{10}$  in both initial and subsequent analyses. The matrix spike/matrix spike duplicate had the internal standard within criteria, although it was on the low side. Since matrix interference was proven both runs for this sample have been reported.

The laboratory followed the USEPA CLP-SOW document #OLMO1.0 for the GC/MS calibration criteria.

<u>Pesticides/PCB's</u> - Forms indicating mass injected for evaluation and individual mixes are enclosed in the package following the Form 10's. MSB SVB-17 12-20' did not meet the required 75 percent recovery for aldrin. Samples SVB-17 12-20', SVB-17 50-60' and corresponding matrix spike/matrix spike duplicates had to be diluted 1:5 because of severe matrix interference. Samples SVB-17 12-20' MS and SVB-17 12-20' MSD had high recovery of several matrix spike compounds because of matrix interference. All calculations were performed off of the DB-1701 column and peaks were taken for calculations of PCB's that are circled on the scans. Corresponding peaks were taken from standards.

The following standards did not meet NYSDEC '89 criteria. After each listed standard, the run was stopped and samples in the last sequence were reanalyzed on the same column.

Date	<u>Time</u>	<u>GC #</u>	<u>Standard</u>	Comments
<b>0</b> 9/19/91	23:46	<b>4</b> A	Ind A	Lindane, heptachlor, aldrin, heptachlor epoxide C <sub>f</sub> >20% dif- ference
09/20/91	00:31	4A	Ind B	delta-BHC C <sub>f</sub> >20% difference

<u>Metals</u> - IEC's are electronically employed by the TJA ICAP-61. However, the ICSA is utilized as a monitoring device to detect any additional adjustments that may be required. These modifications are calculated and applied manually. They are so noted in the raw data.

One "E" flag occurred from serial dilution of sample SVB-17 30-40' for zinc. The diluted sample reading was 48.88 ug/L. This is at the low end of the working curve where instrument drift or electronic fluctuations may be responsible for this flag. It is unlikely that a matrix effect is the cause, however further study would be required to confirm this.

Chromium and copper failed the control limits for duplicate analysis of sample SVB-17 12-20' and resulted in asterisks "\*". This sample contained numerous stones. It is felt that a problem associated with sample homogeneity is responsible for these flags.

Lead and selenium failed the criteria for spike recovery analysis of sample SVB-17 12-20'. The resultant "N" flags are most likely caused by the previously mentioned problems with sample homogeneity. This is substantiated by the fact that the ICAP lead spike recovery for the duplicate sample was low requiring a "W" flag. However, further investigation would be required to prove this.

No other flags or problems were encountered during analysis. All remaining data appears to be consistent.

## **RESULTS**

The results are presented in the following Tables. Also enclosed are the data packages containing all relevant QA/QC and raw data.

Prepared by:

Jeffrey C. Currab Laboratory Manager

JCC/adj

The liability of IEA, Inc. is limited to the actual dollar value of this project.

## TABLE 1.0 30910-1721 UNISYS CORPORATION ANALYTICAL REQUESTS

|--|

## Requested Parameters

SVB-17 20-22', SVB-17 60-62', SVB-17 40-42' 002, SVB-17B, 20-22', SVB-17 40-42' 011, TB 001, TB 002, TB 003, SVB-17 80-82'

SVB-17 12-20', SVB-17 30-40', SVB-17 50-60', SVB-17 70-80'

TCL volatile organics plus a library search for non-target compounds

TCL semi-volatile organics plus a library search for non-target compounds, TCL pesticides/PCB's, TAL metals, cyanide

## TABLE 2.0 30910-1721 UNISYS CORPORATION **EPA TCL VOLATILE ORGANICS**

All values are ug/L.

<u>Dilution Factor</u>	<u> 1.0                                   </u>	
Method Blank I.D.	>G6157 >G6157 >G6157 >G6157	
<u>Compound</u>	Method Quantitation  Method Limits with  Blank TB 001 TB 002 TB 003 Dilution	
Chloromethane Bromomethane Vinyl Chloride Chloroethane Methylene Chloride Acetone Carbon Disulfide 1,1-Dichloroethene 1,1-Dichloroethane 1,2-Dichloroethane 1,2-Dichloroethane 2-Butanone 1,1,1-Trichloroethane Carbon Tetrachloride Vinyl Acetate Bromodichloromethane 1,2-Dichloropropane cis-1,3-Dichloropropene Trichloroethene Dibromochloromethane 1,1,2-Trichloroethane Benzene trans-1,3-Dichloropropene Bromoform 4-Methyl-2-pentanone 2-Hexanone Tetrachloroethene 1,1,2,2-Tetrachloroethane Toluene Chlorobenzene Ethylbenzene	U U U U U 10 0.8J 1JB 2JB 0.9JB 5 6J 2JB U 10 U U U U U 5 U U U U U 5 U U U U U U 5 U U U U	
Styrene Xylene (total)	U U U U 5 U U U U 5	

U, J, B - See Appendix for definition.
 ote: Sample detection limit = quantitation limit x dilution factor.

## TABLE 2.1 30910-1721 UNISYS CORPORATION **EPA TCL VOLATILE ORGANICS**

All values are ug/Kg.

## Sample Identification

<u>Dilution Factor</u>	_1.00	52.00	51.50	_54.50		MEDIUM LEVEL
Method Blank I.D.	>A7470	>A7470	>A7470	>A7_4 <b>7</b> 0		MEDIUM LEVEL
<del></del>				SVB-17		Quantitation
		SVB-17	SVB-17	40-42'		Limits with no
Compound	<u>Blank</u>	20-22′	60-62'	002		<u>Dilution</u>
Chloromethane	U	U	U	U		1,200
Bromomethane	U	U	U	U		1,200
Vinyl Chloride	U	U	U	U	•	1,200
Chloroethane	U	U	U	U		1,200
Methylene Chloride	U	5,700J	U	U		620
Acetone	4,400	130,000B	110,000B	100, <b>0</b> 00B		1,200
Carbon Disulfide	U	U	U	U		620
1,1-Dichloroethene	U	U	U	U		620
1,1-Dichloroethane	U	U	U	U		620
1,2-Dichloroethene (total)	U	U	U	U		620
^hloroform	U	U	U	U		620
,2-Dichloroethane	U	U	U	U		620
2-Butanone	Ų	U	U	U		1,200
1,1,1-Trichloroethane	U	U	U	U		620
Carbon Tetrachloride	Ų	U	U	U		620
Vinyl Acetate	U	U	U	U		1,200
Bromodichloromethane	U	U	U	U		620
1,2-Dichloropropane	U	U	U	U		620
cis-1,3-Dichloropropene	U	U	U	U		620
Trichloroethene	U	92,000	1 <b>3,</b> 000J	Ü		620
Dibromochloromethane	U	U	U	U		620
1,1,2-Trichloroethane	U	U	U	U		620
Benzene	U	U	U	U		620
trans-1,3-Dichloropropene	U	U	U	U		620
Bromoform	U	U	U	U		620
4-methy1-2-Pentanone	U	U	U	U		1,200
2-Hexanone	U	IJ	U	U		1,200
Tetrachloroethene	U	1,200,000		U		620
1,1,2,2-Tetrachloroethane	U	U	U	U		620
Toluene	54J	6,200JB	88,000B	4,600JB		620
Chlorobenzene	U	Ü	U	U		620
Ethylbenzene	U	U	17,000J	U		620
Styrene	U	U	U	U		620
Xylene (total)	U	38,000	98,000	U		620

U, J, B - See Appendix for definition. lote: Sample detection limit = MDL x dilution factor.

## TABLE 2.2 30910-1721 UNISYS CORPORATION **EPA TCL VOLATILE ORGANICS**

All values are ug/Kg.

## Sample Identification

<u>Dilution Factor</u>	1.00	<u>57.00</u>	<u>52.50</u>	55.00		
Mothed Plank I D	> A 7 A 7 O	> A 7 A 7 O	>A7470	>A7470		MEDIUM LEVEL
Method Blank I.D.	<u>&gt;A7470</u>	<u>&gt;A7470</u>	<u>&gt;A/4/U</u>	SVB-17		Quantitation
	Method	SVB-17	SVB-17B	40-42'		Limits with no
<u>Compound</u>		80-82'	20-22'	011		<u>Dilution</u>
<u></u>						
Chloromethane	IJ	U	U	U		1,200
Bromomethane	U	U	U	U		1,200
Vinyl Chloride	U	U	U	U	•	1,200
Chloroethane	U	U	U	U	_	1,200
Methylene Chloride	U	U	5,700J	<b>7</b> ,600J		620
Acetone	4,400	440,000B	130,000B	270,000B		1,200
Carbon Disulfide	Ú	Ú	Ú	Ú		620
1,1-Dichloroethene	U	U	U	U		620
1,1-Dichloroethane	U	U	U	U		620
1,2-Dichloroethene (total)	U	U	U	U		620
îhloroform	U	U	U	U		620
,2-Dichloroethane	U	U	U	U		620
2-Butanone	U	U	U	U		1,200
1,1,1-Trichloroethane	U	U	U	U		620
Carbon Tetrachloride	U	U	U	U		620
Vinyl Acetate	U	U	U	U		1,200
Bromodichloromethane	U	U	U	U		620
1,2-Dichloropropane	U	U	U	U		620
cis-1,3-Dichloropropene	U	U	U	U		620
Trichloroethene	U	U	99,000	U		620
Dibromochloromethane	U	U	U	U		620
1,1,2-Trichloroethane	U	U	U	U		620
Benzene	U	U	U	U		620
trans-1,3-Dichloropropene	U	U	U	U		620
Bromoform	U	U	U	U		620
4-methyl-2-Pentanone	U	U	U	U		1,200
2-Hexanone	U	U	U	U		1,200
Tetrachloroethene	U	U	1,800, <b>0</b> 00E			620
1,1,2,2-Tetrachloroethane	U	U	U	U		620
Toluene	<b>54</b> J	57,000B	5,200JB	U		620
Chlorobenzene	U	U	U	U		620
Ethylbenzene	U	21,000J	U	U		620
Styrene	U	U	U	U		620
Xylene (total)	U	130,000	6,400J	U		620

U, J, B, E - See Appendix for definition.
\_Note: Sample detection limit = MDL x dilution factor.

## TABLE 2.3 30910-1721 UNISYS CORPORATION **EPA TCL VOLATILE ORGANICS**

All values are ug/Kg.

Sample Identification

<u>Dilution Factor</u>	1.00	25.75	57.00	105.00		
						MEDIUM LEVEL
<u>Method Blank I.D.</u>	<u>&gt;A7483</u>	>A7483	>A7483	>A7483		
		SVB-17	SVB-17			Quantitation
		60-62'	80-82'	20-22'		Limits with no
<u>Compound</u>	<u>Blank</u>	<u>RE</u>	<u> </u>	RE		<u>Dilution</u>
Chloromethane	U	U	U	U		1,200
Bromomethane	Ū	Ū	Ü	U		1,200
Vinyl Chloride	U	U	U	U	•	1,200
Chloroethane	ี่ป	U	U	U	-	1,200
Methylene Chloride	U	U	U	U		620
Acetone	2,600	U	U	560,000B		1,200
Carbon Disulfide	Ū	U	U	Ú		620
1,1-Dichloroethene	U	U	U	U		620
1,1-Dichloroethane	U	U	U	U		620
1,2-Dichloroethene (total)	U	4,400J	U	U		620
hloroform	U	Ū	U	U		620
1,2-Dichloroethane	U	U	U	U		620
2-Butanone	U	U	U	U		1,200
1,1,1-Trichloroethane	U	U	U	U		620
Carbon Tetrachloride	U	U	U	U		620
Vinyl Acetate	U	U	U	U		1,200
Bromodichloromethane	U	U	U	U		620
1,2-Dichloropropane	U	U	U	U		620
cis-1,3-Dichloropropene	U	U	U	U		620
Trichloroethene	U	12,000J	U	100,000		620
Dibromochloromethane	U	Ų	U	U		620
1,1,2-Trichloroethane	U	U	U	U		620
Benzene	IJ	U	U	U		620
trans-1,3-Dichloropropene	U	U	U	U		620
Bromoform	Ü	U	U	U		620
4-methyl-2-Pentanone	U	U	U	U		1,200
2-Hexanone	U	U	U	U		1,200
Tetrachloroethene	U	180,000	Ü	2,100,000		620
1,1,2,2-Tetrachloroethane	U	U	U	U 7 5001		620
Toluene	U	82,000	37,000	7,500J		620
Chlorobenzene	U	U 14 0001	U 16 0001	U		620 620
Ethylbenzene	Ų	•	16,000J	Ų		620 620
Styrene	U U	U 06 000	U 000	U 10 0001		620 620
Xylene (total)	U	86,000	93,000	10,000J		620

'', J, B - See Appendix for definition.
ote: Sample detection limit = MDL x dilution factor.

## TABLE 2.4 30910-1721 UNISYS CORPORATION EPA TCL VOLATILE ORGANICS

All values are ug/Kg.

Sample Identification

<u>Dilution Factor</u>	1.00	_1.09_	1.10	1.10	_1.10_		MEDIUM LEVEL
Method Blank I.D.	>A7495	>A7495	>A7495	>A7495	>A7495 SVB-17		
		SVB-17 40-42'					Quantitation Limits with no
Compound					011 MSD		<u>Dilution</u>
Chloromethane	U	U	U	U	U		1,200
Bromomethane	Ŭ	Ŭ	Ŭ	Ŭ	Ŭ		1,200
Vinyl Chloride	Ŭ	Ŭ	Ŭ	Ŭ	Ŭ		1,200
Chloroethane	Ū	Ũ	Ũ	Ü	Ū	•	1,200
Methylene Chloride	Ŭ	300J	340J	Ŭ	Ü	•	620
Acetone	Ŭ	9,900	11,000	5,900	7,600		1,200
Carbon Disulfide	Ū	Ú	Ú	Ü	Ú		620
1,1-Dichloroethene	Ū	Ū	Ū	6,400X	7,300X		620
1,1-Dichloroethane	U	U	U	Ū	Ū		620
1,2-Dichloroethene (total)	U	310J	U	U	U		620
`hloroform	U	U	U	U	U		620
,2-Dichloroethane	U	U	U	U	U		620
2-Butanone	U	U	U	U	U		1,200
1,1,1-Trichloroethane	U	U	U	Ų	U		620
Carbon Tetrachloride	U	U	U	U	U		620
Vinyl Acetate	U	U	U	U	U		1,200
Bromodichloromethane	U	U	U	U	U		620
1,2-Dichloropropane	U	U	U	U	U		620
cis-1,3-Dichloropropene	U	U	U	U	U		620
Trichloroethene	U	490J	U	7,100X	7,900X		620
Dibromochloromethane	U	U	U	Ū	Ū		620
1,1,2-Trichloroethane	U	U	U	U	U		620
Benzene	U	U	U	6,900X	7,900X		620
trans-1,3-Dichloropropene	U	U	U	Ū	Ū		620
Bromoform	U	U	U	U	U		620
4-methy1-2-Pentanone	Ŭ	U	U	U	U		1,200
2-Hexanone	U	U	U	U	U		1,200
Tetrachloroethene	U	3,000	160J	U	U		620
1,1,2,2-Tetrachloroethane	U	U	U	U	U		620
Toluene	U	1,800	U		7,500X		620
Chlorobenzene	U	U	U	6,500X	7,600X		620
Ethylbenzene	U	140J	U	U	U		620
Styrene	U	U	U	U	U		620
Xylene (total)	U	800	U	U	U		620

U, J, X - See Appendix for definition.
\_lote: Sample detection limit = MDL x dilution factor.

## TABLE 3.0 30910-1721 UNISYS CORPORATION VOLATILE TENTATIVELY IDENTIFIED COMPOUNDS

Sample Identification: Method Blank >G6157

CAS#	Compound	<u>RT</u>	Estimated Concentration, ug/L
	None detected		
	Sample Identification:	TB 001	
CAS#	Compound	<u>RT</u>	Estimated <u>Concentration, ug/L</u>
	None detected		-
	Sample Identification:	TB 002	
CAS#	Compound	<u>RT</u>	Estimated Concentration, ug/L
	None detected		
	Sample Identification:	TB 003	
CAS#	Compound	<u>RT</u>	Estimated <u>Concentration, ug/L</u>
	None detected		
	Sample Identification: Metho	od Blank >/	A7470
CAS#	Compound	RT	Estimated <u>Concentration, ug/Kg</u>
	None detected		

## TABLE 3.1 30910-1721 UNISYS CORPORATION VOLATILE TENTATIVELY IDENTIFIED COMPOUNDS

Sample Identification: SVB-17 20-22'

CAS#	Compound		Estimated Concentration, ug/Kg
	Unknown branched alkene	22.33	62,000J
	Unknown branched alkane	23.38	59,000J
	Unknown branched alkene	21.65	45,000J
	Unknown	24.77	45,000J
	Unknown alkene	24.55	44,000J
	Unknown	23.73	44,000J
	Unknown alkane	22.92	43,000J
	Unknown branched alkane	21.42	40,000J
	Unknown cycloalkane	23.83	36,000J
	Unknown C <sub>3</sub> alkyl benzene	22.63	33,000J
	<b>J</b>		

Sample Identification: SVB-17 60-62'

			Estimated
CAS#	Compound	<u>RT</u>	<u>Concentration, ug/Kg</u>
	Unknown alkane	22.90	170,000J
	Unknown C <sub>3</sub> alkyl benzene	23.35	100,000J
	Unknown alkane	24.75	95,000J
	Unknown cycloalkane	21.56	69,000J
	Unknown	22.31	64,000J
	Unknown alkane	20.46	61,000J
	Unknown C, alkyl benzene	22.61	53,000J
	Unknown C <sub>3</sub> alkyl benzene Unknown branched alkane	22.12	50,000J
	Unknown ketone	24.49	44,000J
	Unknown cycloalkane	23.81	40,000J

Sample Identification: SVB-17 40-42' 002

			Estimated	
CAS#	Compound	RT	Concentration, ug/	Κq

None detected

J - See Appendix for definition.

## TABLE 3.2 30910-1721 UNISYS CORPORATION VOLATILE TENTATIVELY IDENTIFIED COMPOUNDS

Sample Identification: SVB-17 80-82'

	Sample Identification:	SVB-17 80-82	•
CAS#	Compound	RT	Estimated <u>Concentration, uq/Kg</u>
	Unknown alkane	22.91	710, <b>0</b> 00J
	Unknown cycloalkane	21.61	280,000J
	Unknown branched alkane	23.40	280,000J
	Unknown alkane	24.80	270,000J
	Unknown alcohol	22.32	250,000J
	Unknown alkane	20.47	230,000J
	Unknown alkane	22.13	220,000J
	Unknown cycloalkane	22.81	200 <del>,</del> 000J
	Unknown cycloalkane	23.85	170+000J
	Unknown branched alkane	21.41	160,000J
	Sample Identification:	SVB-17B 20-22	., Estimated
CAS#	Compound	RT	Concentration, ug/Kg
	oompound		<u>0011001101, 001, 011, 014, 114</u>
	Unknown alkene	24.56	130,000J
	Unknown alkene	23.75	88,000J
	Harley and	00 05	66 6663
	Unknown	23.85	66,000J
	Unknown branched alkene	22.35	65,000J
	Unknown branched alkene Unknown	22.35 24.79	65,000J 65,000J
	Unknown branched alkene Unknown Unknown	22.35 24.79 23.43	65,000J 65,000J 62,000J
	Unknown branched alkene Unknown Unknown Unknown	22.35 24.79 23.43 23.23	65,000J 65,000J 62,000J 56,000J
	Unknown branched alkene Unknown Unknown Unknown Unknown	22.35 24.79 23.43 23.23 24.24	65,000J 65,000J 62,000J 56,000J 56,000J
	Unknown branched alkene Unknown Unknown Unknown Unknown Unknown Unknown	22.35 24.79 23.43 23.23 24.24 21.67	65,000J 65,000J 62,000J 56,000J 49,000J
	Unknown branched alkene Unknown Unknown Unknown Unknown	22.35 24.79 23.43 23.23 24.24	65,000J 65,000J 62,000J 56,000J 56,000J

Sample Identification: SVB-17 40-42' 011

CAS# Compound RT Concentration, ug/Kg

None detected

J - See Appendix for definition.

## TABLE 3.3 30910-1721 UNISYS CORPORATION VOLATILE TENTATIVELY IDENTIFIED COMPOUNDS

Sample Identification: Method Blank >A7483

CAS#	Compound	RT	Estimated <u>Concentration, uq/Kq</u>
	None detected		
	Sample Identification:	SVB-17 60-62'	RE
CAS#	Compound		Estimated <u>Concentration, ug/Kg</u>
	Unknown alkane Unknown C <sub>3</sub> alkyl benzene Unknown alkane Unknown branched alkane Unknown cycloalkane Unknown cycloalkane Unknown C <sub>3</sub> alkyl benzene Unknown C <sub>4</sub> alkyl benzene Unknown alkane Unknown branched alkane	22.94 23.36 20.50 22.32 21.60 22.81 22.61 24.50 24.79 22.16	100,000J 76,000J 47,000J 47,000J 45,000J 44,000J 41,000J 36,000J 36,000J 33,000J

Sample Identification: SVB-17 80-82' RE

CAS#	Compound	RT	Estimated <u>Concentration, ug/Kg</u>
	Unknown alkane	22.90	280,000J
	Unknown alkane	24.75	110,000J
	Unknown	22.31	100,000J
	Unknown cycloalkane	21.56	98,000J
	Unknown alkane	20.46	98,000J
	Unknown branched alkane	22.12	83,000J
		22.61	72,000J
	Unknown C <sub>3</sub> alkyl benzene Unknown branched alkane	21.40	70,000J
	Unknown cycloalkane	22.80	63,000J
	Unknown C <sub>4</sub> alkyl benzene	24.49	57,000J

# TABLE 3.4 30910-1721 UNISYS CORPORATION VOLATILE TENTATIVELY IDENTIFIED COMPOUNDS

Sample Identification: SVB-17B 20-22' RE

CAS#	Compound	<u></u>	Estimated Concentration, ug/Kg
	Unknown alkyl benzene	22.35	620J
	Unknown	23.75	570J
	Unknown ketone Unknown	24.56 24.79	560J 520J
	Olikilowii	24.73	3200
	Sample Identification:	Method Blank >A	\7 <b>4</b> 95
			Estimated
CAS#	Compound	RT	Concentration, ug/Kg
	None detected		
	Sample Identification:	SVB-17 40-42' 0	02 RE
			Estimated
CAS#	Compound	<u>RT</u>	Concentration, ug/Kg
	Unknown alkane	22.93	2,100J
	Unknown alkane	24.79	1,800J
	Unknown branched alkane	23.39	980J
	Unknown alkane	20.49	740J
	Unknown branched alkane	22.31	730J
	Sample Identification:	SVR-17 40-42′ 0	11 RF
	campic racinotification.	010 17 10 12 0	
C N C !!	0	6-	Estimated
<u>CAS#</u>	Compound	<u>RT</u>	Concentration, ug/Kg
	None detected		

J - See Appendix for definition.

#### TABLE 4.0 30910-1741 UNISYS CORPORATION EPA TCL SEMI-VOLATILE ORGANICS

All values are ug/Kg.

Sample Identification

<u>Dilution_Factor</u>	1.00	1.06	1.06	1.09	
Method Blank I.D.	<u>&gt;C1703</u>	<u>&gt;C1703</u>	>C1703	<u>&gt;C1703</u>	
Compound		SVB-17 12-20'		SVB-17 30-40'	Quantitation Limits with no <u>Dilution</u>
Phenol	U	U	U	U	330
bis(2-Chloroethyl)ether	Ŭ	Ŭ	Ŭ	Ŭ	330
2-Chlorophenol	Ŭ	ŭ	Ŭ	Ü	<b>-</b> 330
1,3-Dichlorobenzene	Ŭ	บั	Ŭ	Ŭ	_ 330
1,4-Dichlorobenzene	Ŭ	Ŭ	Ü	Ŭ	330
Benzyl alcohol	Ŭ	Ŭ	Ŭ	Ŭ	330
1,2-Dichlorobenzene	Ŭ	560	550	Ŭ	330
2-Methylphenol	Ü	Ü	Ü	Ū	330
bis(2-Chloroisopropyl)ether	Ū	Ū	Ū	Ū	330
4-Methylphenol	Ŭ	Ŭ	Ū	130J	330
-Nitroso-di-n-propylamine	Ü	Ŭ	Ū	Ü	330
exachloroethane	Ŭ	Ŭ	Ū	Ü	330
Nitrobenzene	Ū	Ū	Ū	U	330
Isophorone	Ū	Ū	U	U	330
2-Nitrophenol	IJ	บ	U	U	330
2,4-Dimethylphenol	Ū	Ū	Ū	Ū	330
Benzoic acid	Ū	Ū	Ū	Ū	1,600
bis(2-Chloroethoxy)methane	Ū	Ū	Ū	Ū	330
2,4-Dichlorophenol	Ū	Ū	Ū	Ü	330
1,2,4-Trichlorobenzene	Ū	Ū	Ū	Ü	330
Naphthalene	Ü	58 <b>0</b>	660	170J	330
4-Chloroaniline	Ü	U	Ü	U	330
Hexachlorobutadiene	Ü	Ü	Ū	U	330
4-Chloro-3-methylphenol	U	U	U	U	330
2-Methylnaphthalene	U	U	430	70J	330
Hexachlorocyclopentadiene	U	Ü	Ŭ	U	330
2,4,6-Trichlorophenol	U	บ	U	U	330
2,4,5-Trichlorophenol	U	U	U	U	1,600
2-Chloronaphthalene	U	U	U	U	330
2-Nitroaniline	U	U	U	U	1,600
Dimethylphthalate	U	U	U	U	330
Acenaphthylene	U	U	U	U	330
2,6-Dinitrotoluene	U	U	U	U	330

U, J - See Appendix for definition. Note: Sample detection limit = quantitation limit x dilution factor.

#### TABLE 4.0 30910-1741 UNISYS CORPORATION **EPA TCL SEMI-VOLATILE ORGANICS**

All values are ug/Kg.

Sample Identification

<u>Dilution Factor</u>	1.00	1.06	1.06	1.09		
Method Blank I.D.	<u>&gt;C1703</u>	<u>&gt;C1703</u>	>C1703	<u>&gt;C1703</u>		
		SVB-17				Quantitation Limits with no
<u>Compound</u>	<u>Blank</u>	<u>12-20'</u>	<u>RE</u>	<u>30-40'</u>		<u>Dilution</u>
3-Nitroaniline	U	U	U	U		1,600
Acenaphthene	U	U	U	U		330
2,4-Dinitrophenol	U	U	U	U	•	1,600
4-Nitrophenol	U	U	U	U	-	1,600
Dibenzofuran	U	U	U	U		330
2,4-Dinitrotoluene	U	U	U	U		330
Diethylphthalate	U	U	U	29J		330
4-Chlorophenyl-phenylether	U	U	U	U		330
Fluorene	U	U	U	U		330
4-Nitroaniline	U	U	U	U		1,600
1,6-Dinitro-2-methylphenol	U	U	U	U		1,600
Nitrosodiphenylamine (1)	U	U	U	U		330
4-Bromophenyl-phenylether	U	U	U	U		330
Hexachlorobenzene	U	U	U	U		330
Pentachlorophenol	Ú	U	U	U		1,600
Phenanthrene	Ü	Ü	260J	61J		<sup>2</sup> 330
Anthracene	Ü	Ü	Ü	U		330
Di-n-butylphthalate	13J	270JB	270JB	<b>60</b> JB		330
Fluoranthene	Ü	Ü	U	U		330
Pyrene	Ü	Ü	Ŭ	Ü		330
Butylbenzylphthalate	Ŭ	Ŭ	Ü	Ŭ		330
3,3'-Dichlorobenzidine	Ŭ	Ŭ	Ŭ	Ū		660
Benzo(a) anthracene	Ŭ	Ŭ	Ŭ	Ŭ		330
Chrysene	Ŭ	ŭ	Ŭ	Ŭ		330
bis(2-Ethylhexyl)phthalate	220J	2.600B	2,600B	640B		330
Di-n-octylphthalate	U	U	U	U		330
Benzo(b)fluoranthene	Ŭ	Ŭ	Ŭ	Ŭ		330
Benzo(k)fluoranthene	Ŭ	Ŭ	Ŭ	Ŭ		330
Benzo(a)pyrene	Ŭ	Ŭ	Ŭ	Ŭ		330
Indeno(1,2,3-cd)pyrene	Ŭ	Ŭ	Ŭ	Ŭ		330
Dibenzo(a,h)anthracene	Ŭ	Ŭ	Ŭ	Ŭ		330
Benzo(g,h,i)perylene	Ŭ	Ŭ	Ŭ	Ŭ		330

U, J, B, (1) - See Appendix for definition.

Note: Sample detection limit = quantitation limit x dilution factor.

#### TABLE 4.1 30910-1721 UNISYS CORPORATION **EPA TCL SEMI-VOLATILE ORGANICS**

All values are ug/Kg.

#### Sample Identification

<u>Dilution Factor</u>	1.00	1.06	1.08	1.06	1.06		
Method Blank I.D.	<u>&gt;C1703</u>	<u>&gt;C1703</u>	>C1703	<u>&gt;C1703</u>	>C1703		
<u>Compound</u>		SVB-27 50-60'		SVB-17 12-20' <u>MS</u>	SVB-17 12-20' <u>MSD</u>		Quantitation Limits with no <u>Dilution</u>
Phenol	U	U	U	3,800X	4,400X		330
bis(2-Chloroethyl)ether	U	U	U	Ú	Ū		330
2-Chlorophenol	U	U	U	3,600X	4,100X	•	330
1,3-Dichlorobenzene	U	U	U	Ū	Ū	_	330
1,4-Dichlorobenzene	Ų	U	U	1,400X	1,700X		330
Benzyl alcohol	U	U	U	Ū	Ū		330
1,2-Dichlorobenzene	U	1,100	620	500	590		330
2-Methylphenol	U	U	U	U	U		330
bis(2-Chloroisopropyl)ether	U	U	U	IJ	U		330
4-Methylphenol	U	U	U	U	U		330
'-Nitroso-di-n-propylamine	U	U	U	2,800X	2,700X		330
lexachloroethane	U	U	U	Ū	U		330
Nitrobenzene	U	U	U	U	U		330
Isophorone	U	U	U	U	U		330
2-Nitrophenol	U	U	U	U	U		330
2,4-Dimethylphenol	IJ	310J	U	U	U		330
Benzoic acid	U	U	U	U	U		1,600
bis(2-Chloroethoxy)methane	U	U	U	U	U		330
2,4-Dichlorophenol	U	U	U	U	U		330
1,2,4-Trichlorobenzene	U	U	U	2,300X	2,600X		330
Naphthalene	Ü	3,100	2,600	<b>5</b> 00	530		330
4-Chloroaniline	U	Ú	Ú	U	U		330
Hexachlorobutadiene	U	U	U	U	U		330
4-Chloro-3-methylphenol	U	U	U	4,000X	4,000X		330
2-Methylnaphthalene	U	1,500	1,600	330J	320J		330
Hexachlorocyclopentadiene	U	Ú	Ū	U	U		330
2,4,6-Trichlorophenol	U	U	Ü	U	U		330
2,4,5-Trichlorophenol	U	U	IJ	U	U		1,600
2-Chloronaphthalene	U	U	U	U	U		330
2-Nitroaniline	U	U	U	U	U		1,600
Dimethylphthalate	U	U	U	U	U		330
Acenaphthylene	U	U	U	U	U		330
2,6-Dinitrotoluene	IJ	U	U	U	U		330

U, J, X - See Appendix for definition. Note: Sample detection limit = quantitation limit x dilution factor.

#### TABLE 4.1 30910-1721 UNISYS CORPORATION EPA TCL SEMI-VOLATILE ORGANICS

All values are ug/Kg.

Sample Identification

<u>Dilution Factor</u>	1.00	1.06	1.08	1.06	1.06		
Method Blank I.D.	>C1703	>C1703	>C1703	>C1703	>C1703_		
	Mathad	CVD 07	CVD 17	SVB-17	SVB-17 12-20'		Quantitation
<u>Compound</u>		SVB-27 50-60'		12-20' MS	MSD		Limits with no Dilution
<u></u>	2. 41,11	<u>• • • • • • • • • • • • • • • • • • • </u>					
3-Nitroaniline	U	U	U	U	U		1,600
Acenaphthene	U	U	U	2,700X	2,500X		330
2,4-Dinitrophenol	U	U	U	U	U	•	1,600
4-Nitrophenol	U	U	U	9,100XE	8,300XE	-	1,600
Dibenzofuran	U	U	Ų	U	U		330
2,4-Dinitrotoluene	IJ	U	U	2,000X	1,900X		330
Diethylphthalate	U	U	U	U	U		330
4-Chlorophenyl-phenylether	U	U	U	U	U		330
Fluorene	U	110J	150J	U	U		330
<pre>-Nitroaniline</pre>	U	U	U	U	U		1,600
,6-Dinitro-2-methylphenol	U	U	IJ	U	U		1,600
N-Nitrosodiphenylamine (1)	U	U	U	U	U		330
<pre>4-Bromophenyl-phenylether</pre>	U	U	U	U	U		330
Hexachlorobenzene	U	U	U	U	U		330
Pentachlorophenol	U	U	Ü		6, <b>4</b> 00XE		1,600
Phenanthrene	U	570	550	230J	U		330
Anthracene	U	U	U	U	U		330
Di-n-butylphthalate	13J	350B	170JB	280JB	230JB		330
Fluoranthene	U	U	U	U	Ü		330
Pyrene	U	U	U	1,600X	1,700X		330
Butylbenzylphthalate	U	U	U	220J	220J		330
3,3'-Dichlorobenzidine	U	U	U	U	U		660
Benzo(a)anthracene	U	U	U	U	U		330
Chrysène	U	U	U	U	U		330
bis(2-Ethylhexyl)phthalate	220J	2,200B	<b>520</b> B	2,300B	2,900B		330
Di-n-octylphthalate	U	U	U	U	U		330
Benzo(b)fluoranthene	U	U	U	U	U		330
Benzo(k)fluoranthene	U	U	U	U	U		330
Benzo(a)pyrene	U	U	U	U	U		330
Indeno(1,2,3-cd)pyrene	U	U	U	U	U		330
Dibenzo(a,h)anthracene	U	U	U	U	U		330
Benzo(g,h,i)perylene	U	U	U	U	U		330

U, J, B, X, E, (1) - See Appendix for definition. Note: Sample detection limit = quantitation limit x dilution factor.

## TABLE 5.0 30910-1721 UNISYS CORPORATION SEMI-VOLATILE TENTATIVELY IDENTIFIED COMPOUNDS

Sample Identification: Method Blank >C1703

CAS#	Compound	RT	Estimated <u>Concentration, ug/Kg</u>		
	Aldol condensation product	8.82	33,000JA		
	Unknown acid ester	30.10	300J		
	Unknown ketone (MW=98)	7.62	190J		

Sample Identification: SVB-17 12-20'

			Estimated
CAS#	Compound	<u>RT</u>	Concentration, ug/Kg
	Unknown alkane	22.91	50,000J
		19.08	•
	Unknown	_	14,000J
	Unknown	19.70	12,000J
	Unknown alkane	18.91	11,000J
	Aldol condensation product	8.89	6,200JAB
	Unknown	20.23	5,000J
	Unknown alkane	20.30	4,800J
	Unknown alkane	21.63	4,600J
	Unknown branched alkane	22.27	3,800J
	Unknown branched alkane	19.78	3,000J
	Unknown C <sub>4</sub> alkyl benzene	14.56	2,700J
	Unknown branched alkane	18.59	1,900J
	Unknown	12.97	1,900J
	Unknown C <sub>3</sub> alkyl benzene	12.21	1,700J
	Unknown	14.86	1,500J
	Unknown	18.21	1,500J
	Unknown alkane	14.13	1,500J
	Unknown branched alkene	11.43	1,500J
	Unknown alkane	17.41	1,400J
	Unknown	20.40	1,100J
	Unknown	14.41	1,100J

### TABLE 5.1 30910-1721 UNISYS CORPORATION SEMI-VOLATILE TENTATIVELY IDENTIFIED COMPOUNDS

Sample Identification: SVB-17 12-20' RE

CAS#	Compound	RT	Estimated <u>Concentration, ug/Kg</u>
	Unknown alkane	18.94	11,000J
	Unknown	19.72	9,700J
	Unknown	19.09	6,600J
	Unknown alkene	20.24	4,900J
	Aldol condensation product	8.87	4,600JAB
	Unknown branched alkane	21.66	4,500J
	Unknown	23.15	4,300J
	Unknown alkane	20.32	4, <del>2</del> 00J
	Unknown branched alkane	22.87	3, <b>3</b> 00J
	Unknown branched alkane	22.30	3,100J
	Unknown	14.58	2,700J
	Unknown branched alkane	19.81	2,600J
	Unknown branched alkane	18.62	1,900J
	Unknown branched alkane	13.01	1,700J
	Unknown	14.88	1,600J
	Unknown alkane	12.23	1,500J
	Unknown branched cycloalkane	18.24	1,400J
	Unknown alkane	17.44	1,400J
	Unknown branched alkene	11.45	1,300J
	Unknown branched alkane	18.40	1,200J
	Unknown alkane	14.15	1,200J

Sample Identification: SVB-17 30-40'

			Estimated
CAS#	Compound	<u>RT</u>	Concentration, uq/Kq
	Aldol condensation product	8.90	47,000JAB
1120214	Undecane	22.95	<b>3</b> ,600J
	Unknown alkane	14.05	2,700J
	Unknown alkane	12.17	2,400J
	Unknown	19.09	2,100J
	Unknown branched alkane	22.88	1,600J
629787	Heptadecane	22.80	1,600J
	Unknown branched alkane	23.99	1,600J
	Unknown	18.99	1,500J
	Unknown alkane	25.10	1,400J
	Unknown alkane	21.55	1,100J
	Unknown alkane	24.11	1,000J
	Unknown	19.64	930J

### TABLE 5.2 30910-1721 UNISYS CORPORATION SEMI-VOLATILE TENTATIVELY IDENTIFIED COMPOUNDS

Sample Identification: SVB-17 30-40' (Continued)

Estimated <u>Concentration, ug/Kg</u>
8 890J
3 840J
2 790J
5 730J
5 700J
5 670J
4 650J
9 620J

Sample Identification: SVB-17 50-60'

			Estimated
CAS#	<u>Compound</u>	RT	Concentration, ug/Kg
	Unknown alkane	22.88	34,000J
	Unknown branched alkane	24.07	16,000J
	Unknown alkane	25.18	12,000J
	Unknown branched alkane	24.19	9,500J
	Unknown alkane	18.88	8,200J
	Unknown alkane	21.60	8,100J
	Unknown	19.05	8,000J
	Unknown alkane	26.24	7,700J
	Unknown	19.69	7,600J
	Aldol condensation product	8.84	5,900JAB
	Unknown alkane	20.29	5,300J
	Unknown alkane	12.25	4,400J
	Unknown alkane	27.23	3,700J
	Unknown branched alkane	23.98	3,600J
	Unknown	20.20	3,600J
	Unknown C <sub>4</sub> alkyl benzene	13.32	2,900J
	Unknown branched alkane	22.24	2,800J
	Unknow <b>n</b>	11.49	2,700J
	Unknown alkane	14.12	2,700J
	Unknown PAH	19.57	2,400J
	Unknown alkane	19.75	2,300J

# TABLE 5.3 30910-1721 UNISYS CORPORATION SEMI-VOLATILE TENTATIVELY IDENTIFIED COMPOUNDS

Sample Identification: SVB-17 70-80'

CAS#	Compound	RT	Estimated Concentration, ug/Kg
	Unknown branched alkane	22.88	63, <b>0</b> 00J
	Unknown branched alkane	24.05	38,000J
	Aldol condensation product	8.92	22,000JAB
	Unknown branched alkane	22.94	21,000J
	Unknown alkane	25.17	19,000J
	Unknown alkane	24.18	15,00 <b>0</b> J
	Unknown alkane	21.60	13,000J
	Unknown alkane	26.23	11 <b>,</b> 000J
	Unknown branched alkane	12.23	9, <b>4</b> 00J
	Unknown alkane	14.12	7,900J
	Unknown alkane	25.09	7,900J
	Unknown alkane	18.88	7,300J
	Unknown branched alkane	20.28	5,500J
	Unknown C∠ alkyl benzene	13.31	5,200J
	Unknown branched alkane	22.22	5,000J
	Unknown C <sub>3</sub> alkyl benzene	11.49	4,600J
	Unknown	20.20	3,900J
	Unknown	19.67	3,600J
	Unknown	19.10	3,400J
	Unknown branched alkane	13.44	3,300J
	Unknown branched alkane	12.66	3,100J

#### TABLE 6.0 30910-1721 UNISYS CORPORATION EPA TCL PESTICIDES/PCB's

All values are ug/Kg.

#### Sample Identification

<u>Dilution Factor</u>	1.00	5.30	1.09	5.30	1.08	5.30	5.30		
Method Blank I.D.	0827 -B02	0827 -B02	0827 -B02	0827 -B02	0827 B02	0827 -B02	0827 <u>-B02</u>		
	Met hod	SVB-17	SVR-17	SVR-18	SVR-17	SVB-17 12-20'	SVB-17 12-20'		Quantitation Limits with no
<u>Compound</u>		12-20'				MS	MSD		Dilution
alpha-BHC	U	U	U	U	U	U	U		8.0
beta-BHC	U	U	U	U	U	U	U		8.0
delta-BHC	U	U	U	U	Ü	U	U	-	8.0
gamma-BHC	U	48	U	88	U	98X	110X	_	8.0
Heptachlor	U	U	U	U	U	64X	68X		8.0
Aldrin	U	U	U	U	U	98X	110X		8.0
Heptachlor Epoxide	U	U	U	U	U	U	U		8.0
Endosulfan I	U	U	U	U	U	U	U		8.0
Dieldrin	U	U	U	U	U	74X	81X		16
,4′-DDE	U	U	U	U	U	U	U		16
<b>√</b> ndrin	U	U	U	U	U	99X	100X		16
Endosulfan II	U	U	U	U	U	IJ	U		16
4,4'-DDD	U	U	U	U	U	U	U		16
Endosulfan Sulfate	U	U	U	U	U	U	U		16
4,4′-DDT	U	U	U	U	U	240X	240X		16
Methoxychlor	U	U	U	U	U	U	U		80
Endrin-Ketone	U	U	U	IJ	U	U	U		16
alpha-Chlordane	U	U	U	U	U	U	U		80
gamma-Chlordane	U	U	U	U	U	U	U		80
Ťoxaphene	U	U	U	U	U	U	U		160
PCB - 1016	U	U	U	U	U	U	U		80
PCB - 1221	U	U	U	U	U	U	U		80
PCB - 1232	U	U	U	U	U	U	U		80
PCB - 1242	U	U	U	U	U	U	U		80
PCB - 1248	U	U	U	2,300	2,000	U	U		80
PCB - 1254	U	890	U	760	730	1,000	1,100		160
PCB - 1260	. <b>U</b>	U	U	U	U	Ū	Ū		160

U, X - See Appendix for definition. Note: Sample detection limit = quantitation limit x dilution factor.

#### TABLE 7.0 30910-1721 UNISYS CORPORATION TAL METALS PLUS CYANIDE

All values are mg/Kg dry basis.

<u>Parameter</u>	SVB-17 12-20'	SVB-17 30-40'	SVB-17 50-60'	SVB-17_70-80'
Aluminum	2,100	5,430	2,120	1,750
Antimony	3.9U	4.0U	4.1U	3.9U
Arsenic	0.32B	0.57B	0.18U	<b>0</b> .18U
Barium	13.5B	37.5	18.7B	13.5B
Beryllium	0.18U	0.418	0.19B	0.18U
Cadmium	0.58B	0.52B	0.19U	~0.18U
Calcium	393B	961	196B	104B
Chromium	8.7*	13.2*	4.1*	5.1*
Cobalt	4.1B	5.5B	1.6B	1.7B
Copper	24.1*	15.6*	5.1*	3.8B*
Iron	7,430	12,800	4,570	4,910
Lead	5.3N	22.3N	2.2N	2.7NS
Magnesium	812B	1,900	566B	<b>549</b> B
Manganese	61.9	218	93.1	58.4
Mercury	0.09	0.09U	0.09U	<b>0.08</b> U
Nickel	18.4	15.0	7.9	8.3
Potassium	444B	1,370	459B	392B
Selenium	0.18UN	0.18UNW	0.18UN	0.18UN
Silver	0.54U	<b>0.55</b> U	0.56U	0.54U
Sodium	33.8B	96.7B	<b>38.7</b> B	62.9B
Thallium	0.35UW	0.37UW	0.36UW	0.36UW
Vanadium	6.3B	16.4	6.1B	5.5B
Zinc	27.0E	32.3E	10.8E	9.1E
Cyanide	2.6U	2.7U	2.7U	2.6U

B, E, N, S, U, W, \* - See Metals Appendix for definition.

#### **APPENDIX**

- U Indicates that the compound was analyzed for but not detected.
- J Indicates that the compound was analyzed for and determined to be present in the sample. The mass spectrum of the compound meets the identification criteria of the method. The concentration listed is an estimated value, which is less than the specified minimum detection limit but is greater than zero.
- B This flag is used when the analyte is found in the blanks as well as the sample. It indicates possible sample contamination and warns the data user to use caution when applying the results of this analyte.
- N Indicates that the compound was analyzed for but not requested as an analyte. Value will not be listed on tabular result sheet.
- X Matrix spike compound.
- (1) Cannot be separated.
- (2) Decomposes to azobenzene. Measured and calibrated as azobenzene.
- A This flag indicates that a TIC is a suspected aldol condensation product.
- E Indicates that it exceeds calibration curve range.
- D This flag identifies all compounds identified in an analysis at a secondary dilution factor.
- C Confirmed by GC/MS.
- T Compound present in TCLP blank.

#### APPENDIX/METALS DATA

#### C - Concentration qualifiers

- U Indicates analyte result less than instrument detection limit (IDL)
- B Indicates analyte result between IDL and contract required detection limit (CRDL)

#### Q - QC qualifiers

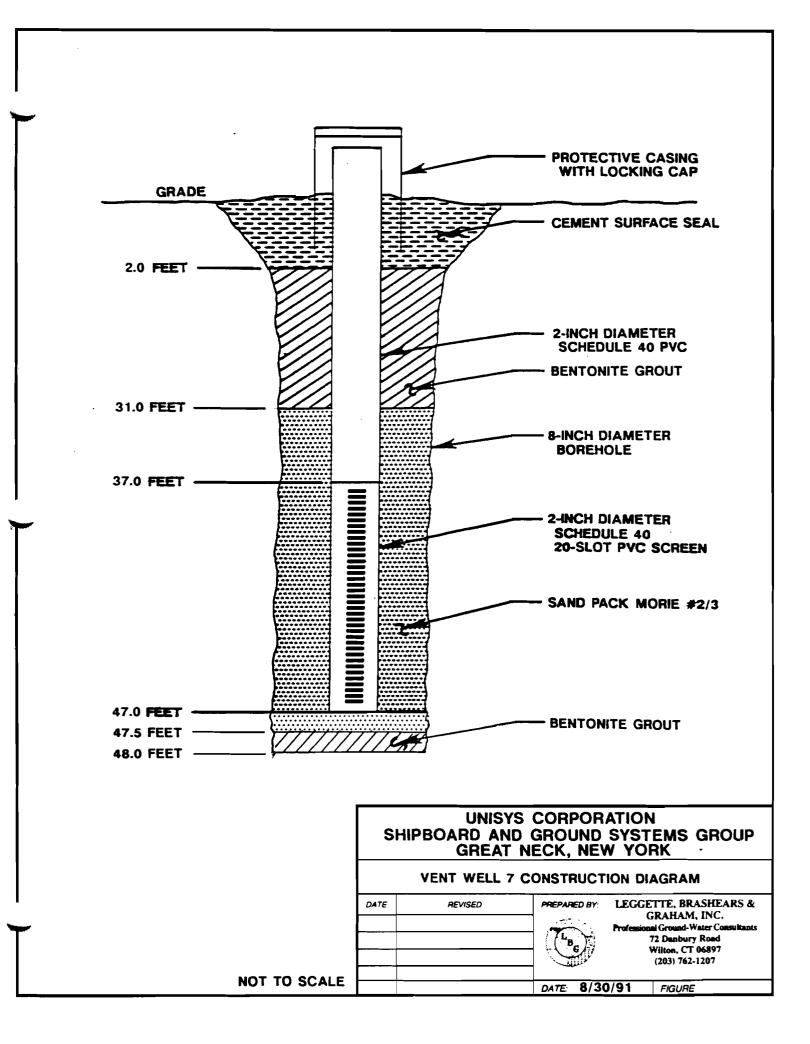
- E Reported value is estimated because of the presence of interference
- M Duplicate injection precision not met
- N Spiked sample recovery not within control limits
- S The reported value was determined by the method of standard additions (MSA)
- W Post-digest spike recovery furnace analysis was out of 85-115 percent control limit, while sample absorbance was less than 50 percent of spike absorbance
- \* Duplicate analysis not within control limit
- + Correlation coefficient for MSA is less than 0.995

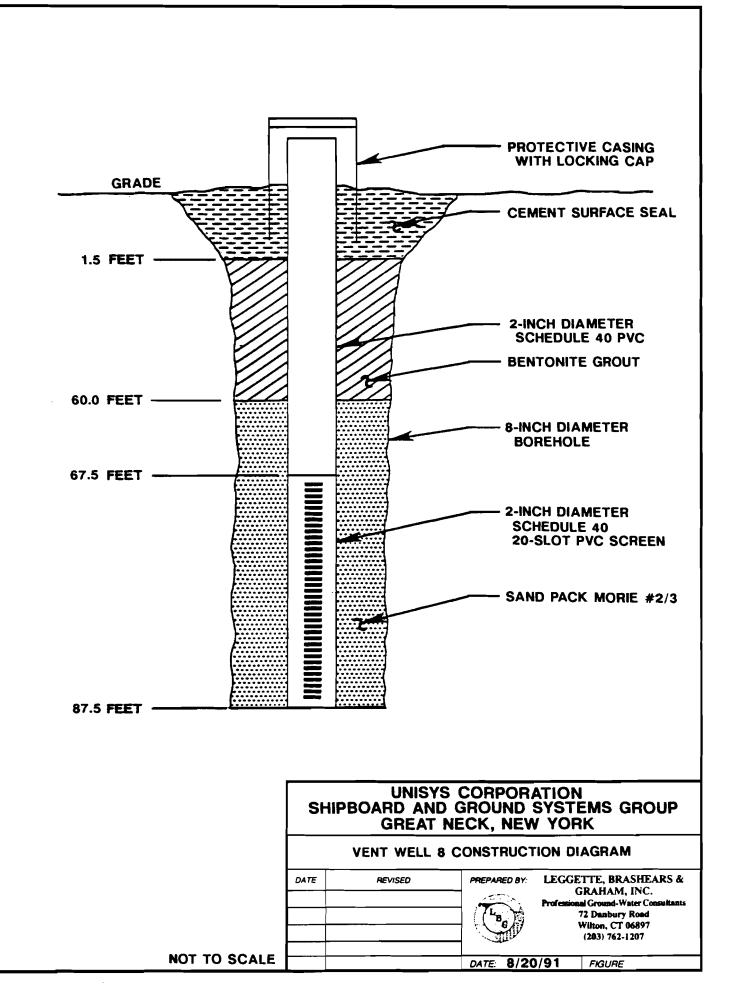
#### M - Method codes

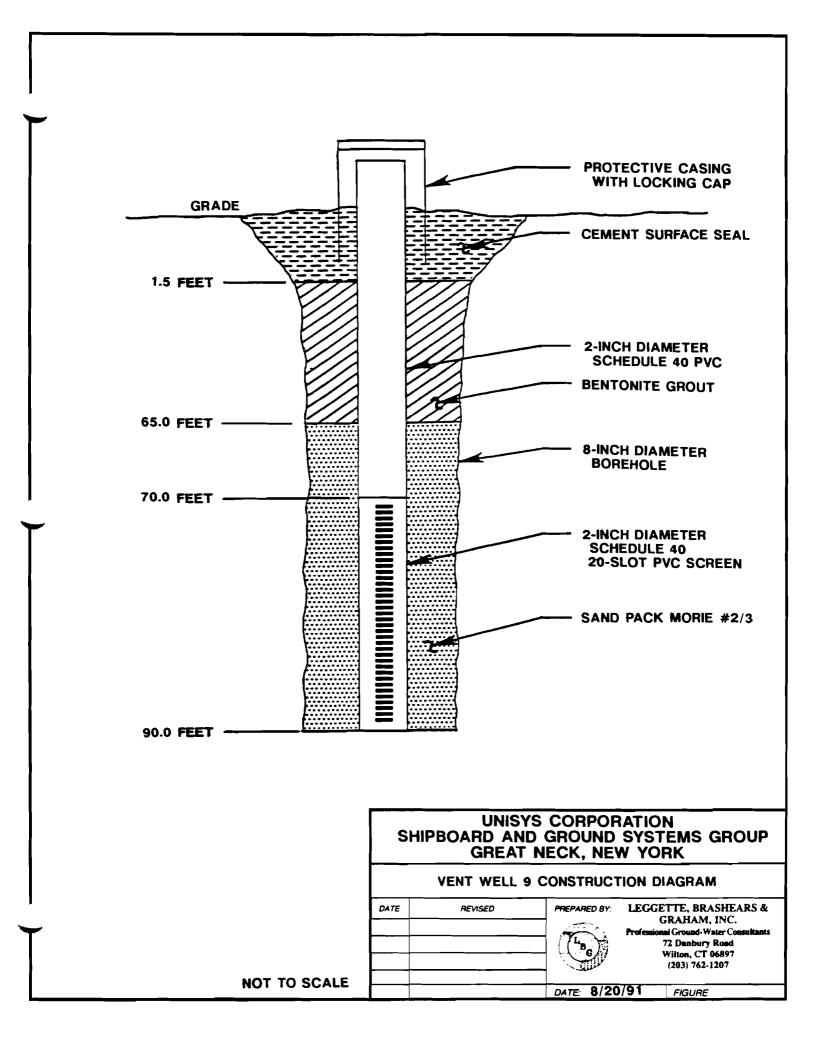
- P ICP
- A Flame AA
- F Furnace AA
- CV Cold vapor AA (manual)
  - C Cyanide
- NR Not Required
- NC Not Calculated as per protocols

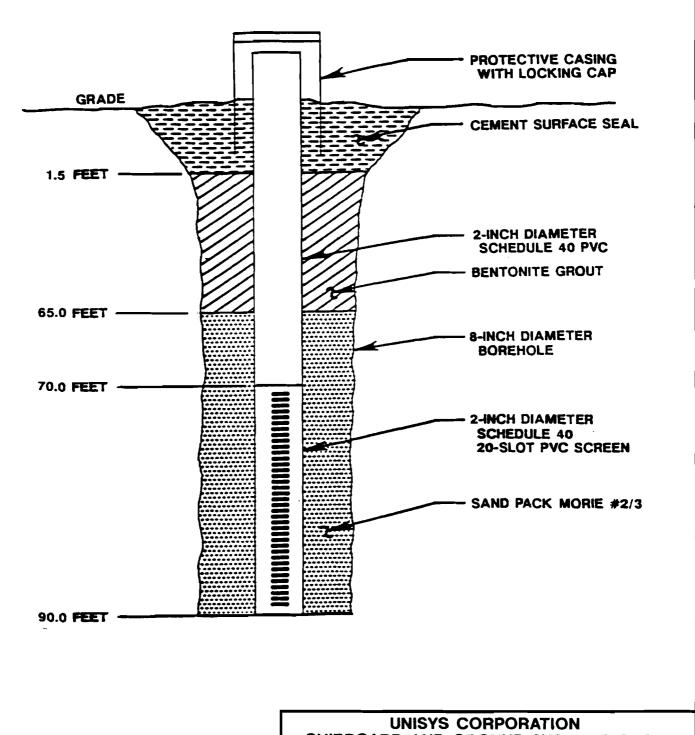
#### APPENDIX L3

# LBG 1991 VENT WELL CONSTRUCTION LOGS









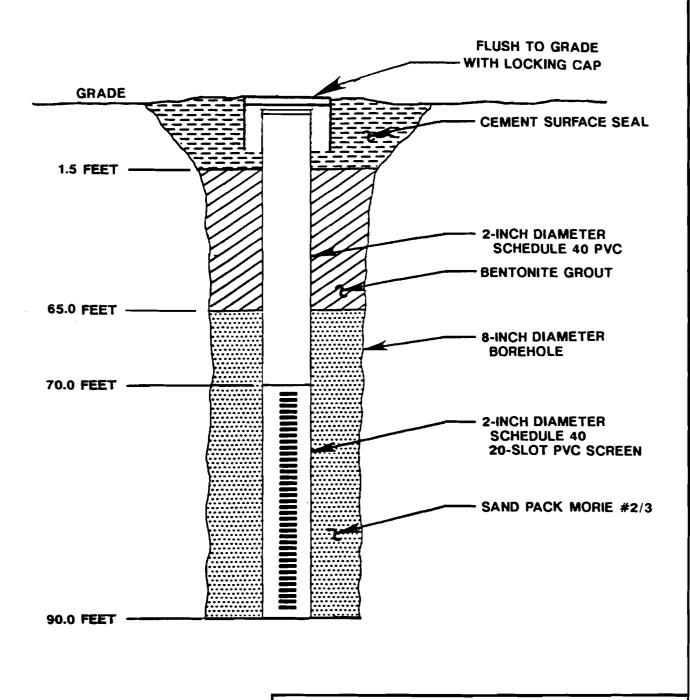
#### **VENT WELL 10 CONSTRUCTION DIAGRAM**

DATE REVISED PREPARED BY:

LEGGETTE, BRASHEARS & GRAHAM, INC.
Professional Ground-Water Consultants
72 Danbury Road
Wilton, CT 06897
(203) 762-1207

NOT TO SCALE

DATE: 8/27/91 FIGURE



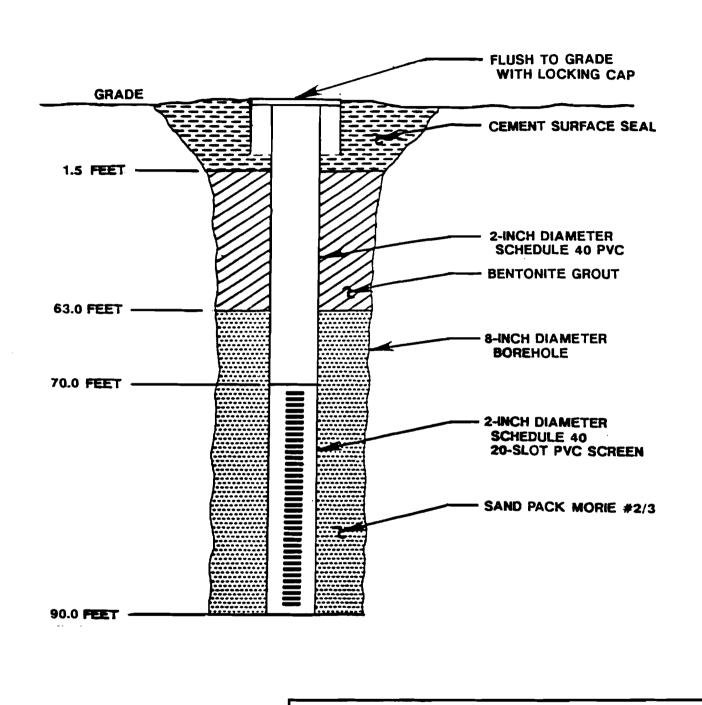
#### **VENT WELL 11 CONSTRUCTION DIAGRAM**

DATE REVISED

PREPARED BY: LEGGETTE, BRASHEARS & GRAHAM, INC.
Professional Ground-Water Consultants
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Wilton, CT 06897
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DATE: 8/20/91 FIGURE

NOT TO SCALE



#### **VENT WELL 12 CONSTRUCTION DIAGRAM**

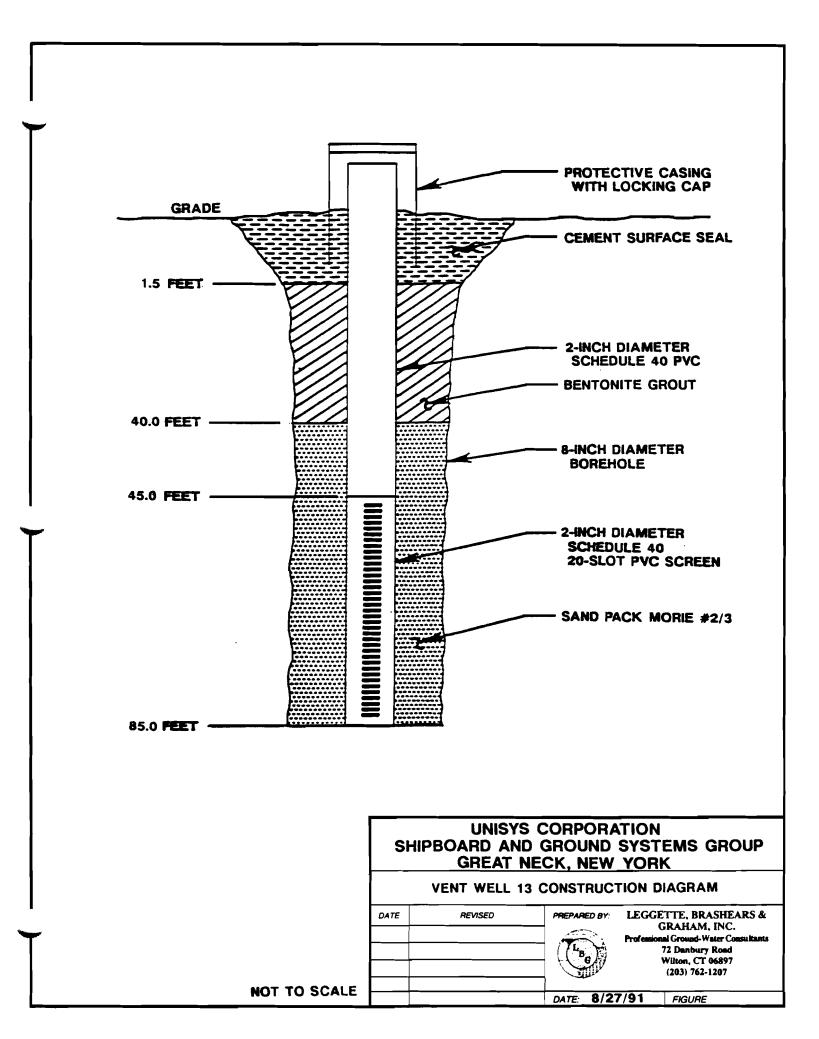
DATE REVISED

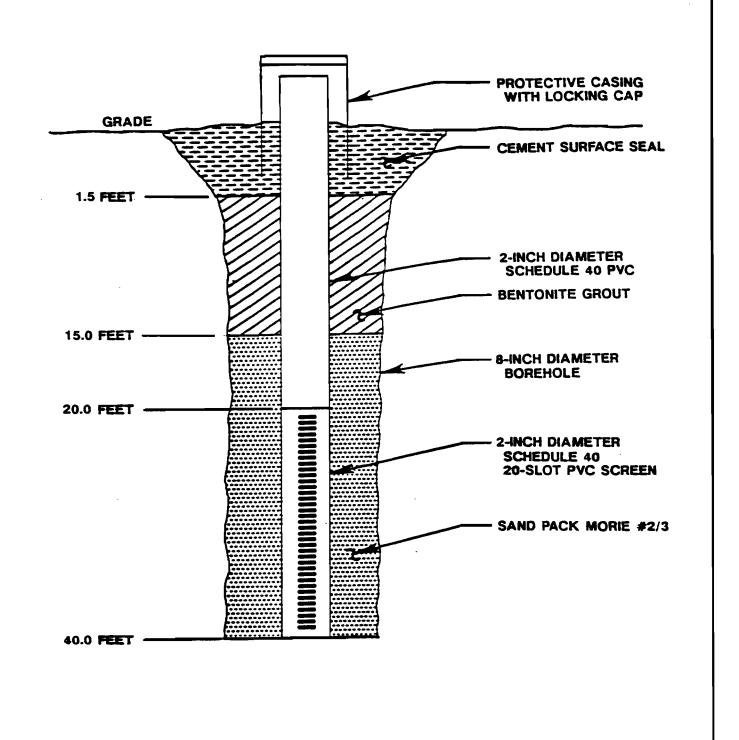
PREPARED BY: LEGGETTE, BRASHEARS & GRAHAM, INC.

Professional Ground-Water Consultants
72 Danbury Road
Wilton, CT 06897
(203) 762-1207

DATE: 8/27/91 FIGURE

NOT TO SCALE





#### **VENT WELL 14 CONSTRUCTION DIAGRAM**

DATE REVISED PREPARED BY: LEGGETTE, BRASHEARS & GRAHAM, INC. onal Ground-Water Consultants

**NOT TO SCALE** 

DATE: 8/27/91 FIGURE

72 Danbury Road Wilton, CT 06897 (203) 762-1207

# APPENDIX L4 LBG 1991 WELL LOGS

#### **GEOLOGIC LOG**

LEGGETTE, BRASHEARS & GRAHAM, INC.

OWNER Unisys Corporation, Great Neck, NY

**PAGES** 

OF 2

WELL NO. 2 MU

PAGE

WILTON, CONNECTICUT

SCREEN TYPE PVC Schedule-40 LOCATION Northeast corner of the

SLOT NO. 20 DIAM. garage building 4-inch

DATE COMPLETED July 11, 1991 SETTING 175-185 feet below grade

DRILLING SAND PACK 2-3 Morie Mix; 168-185 ft bg R & L Well Drilling

COMPANY CASING 4-inch Schedule-40 PVC

DRILLING

METHOD Mud rotary/8-inch annulus SETTING 0 - 175 feet below grade

SAMPLING

METHOD Grab/wash samples DEVELOPMENT July 12, 1991; submersible pump

OBSERVER Keith Yocis DURATION

REFERENCE POINT (RP) STATIC WATER LEVEL

Grade ELEVATION OF RP YIELD Greater than 20 gpm

Bentonite grout set from 0 to 168 feet below grade. Drilling fluid; REMARKS Aquagel Gold Seal, bentonite mud.

DEPTH (FEET	ТО	DESCRIPTION
	10	Gravel, fine to coarse, and medium to very coarse sand; little very
		fine to fine sand; trace silt; tan.
	20	GRAVEL, fine, little medium, and medium to very coarse sand; trace
		very fine to fine sand and silt; tan.
	30	GRAVEL, fine, little medium gravel, and medium to very coarse sand;
		trace very fine to fine sand and silt; tan.
	40	Gravel, fine, and coarse to very coarse sand, little fine to medium;
		trace very fine sand and silt; tan.
	50	Gravel, fine, trace medium; very coarse to coarse sand, little fine to
_		medium; trace very fine sand and silt; clay; tan.

OWNER	Unisys	Corporation
WELL NO.	2 MU	PAGE 2 OF 2 PAGES
EPTH (FE	ET) TO	DESCRIPTION
	70	Gravel, fine, and very coarse sand, little medium; trace very fine
		to fine sand and silt; tan.
	80	Gravel, fine, and very coarse sand, little medium; trace very fine
		to fine sand and silt; MICAS; tan.
	90	GRAVEL, fine, some medium; some very coarse sand, little very fine;
		trace silt; MICAS; tan.
	100	Gravel, fine, and coarse sand; little very fine to medium sand and
_		sandy clay; trace MICAS; tan.
	120	SAND, fine to medium, trace coarse; MICAS; tan.
-	130	SAND, fine to medium, trace coarse; brown.
	140	SAND, fine to medium, trace coarse; MICAS; tan.
	150	SAND, fine to medium, trace coarse; MICAS; tan.
,——	160	SAND, fine to medium, trace coarse; MICAS; tan.
	170	SAND, fine to medium, little coarse, trace very coarse; MICAS; tan.
176	179	CLAY; trace lignite; dark gray.
	180	SAND, very fine to fine; trace silt; MICAS; tan.
	190	SAND, very fine to fine; trace silt; MICAS; tan.
	195	SAND, very fine, little fine; trace silt; and MICAS; tan.
	200	SILT.
	205	SAND, medium, little fine to coarse; trace fine gravel; tan.
	210	SAND, medium to coarse, some fine, little very coarse; trace fine
		gravel; brown.
	210	End of boring.
<del></del>		

# **GEOLOGIC LOG** LEGGETTE, BRASHEARS & GRAHAM, INC. WELL NO. 6 MI

OWNER Unisys Corporation
Great Neck, New York

	6 M1			
WILTON, CONNECTICUT	PAGE 1 OF 2 PAGES			
LOCATION West of the former foundry building, 30 feet north of	SCREEN TYPE PVC Schedule 40			
diffusion Well 5	DIAM. 4-inch SLOT NO. 20			
DATE COMPLETED June 28, 1991	SETTING 215-235 feet below grade			
DRILLING R & L Well Drilling	SAND PACK 175-240 feet below grade			
DRILLING	CASING 4-inch PVC Schedule 40			
METHOD Mud rotary/8-inch annulus	2.5 feet above grade to			
SAMPLING	SETTING 215 feet below grade			
METHOD Grab/wash samples	DEVELOPMENT June 28, 1991 submersible pump			
OBSERVER J. Benvegna	DURATION 1.25 hours			
REFERENCE POINT (RP) Grade	STATIC WATER LEVEL 80.4 feet below grade			
ELEVATION OF RP	YIELD Greater than 20 gallons per minute			

Bentonite grout set from 0 - 175 feet below grade. Drilling fluid: aquagel gold REMARKS seal bentonite mud.

DEPTH (FEE FROM	T)   TO	DESCRIPTION
	10	SAND, coarse to very coarse; some fine gravel.
	20	SAND, coarse; some very coarse; little fine gravel.
	25	SAND, coarse; little very coarse; some silt; little clay; brown.
	30	SAND, coarse; some very coarse; little medium; trace silt.
	40	SAND, coarse to very coarse; some medium; little fine gravel; trace
<del>-</del>	-	silt and clay.
	40	Clean mud pit, noted; fine to very fine sand not in grab samples.
	50	SAND, coarse, some medium to very coarse; trace gravel.
	60	SAND, coarse to very coarse, little medium; trace fine gravel.
	70	SAND, coarse, some medium, trace very coarse; slightly finer than
		60 foot sample.
	80	SAND, coarse, some medium, trace very coarse.

OWNER	Unisys	Corporation
WELL NO.	6 MI	PAGE 2 OF 2 PAGES
DEPTH (FEET)		DESCRIPTION
	90	SAND, medium, some coarse, little fine.
	100	SAND, fine to medium, some very fine; very little sample, return;
		washing through strainer.
	110	SAND, fine to very fine; brown.
	120	SAND, fine to very fine, trace medium; brown.
	130	SAND, very fine to fine; tan, gray.
	140	SAND, very fine to fine; tan, gray.
	150	SAND, very fine to fine; gray.
	170	SAND, very fine to fine; gray.
	190	SAND, fine to very fine, little medium; tan; slightly coarser.
	200	SAND, fine, some very fine to medium; tan.
	210	SAND, fine, some very fine to medium; tan.
	220	SAND, medium, some fine; gray. Visually coarser.
	230	SAND, medium, trace fine; gray.
	240	SAND, coarse, little very coarse, trace medium.
	240	End of boring.
-	<del></del>	
- <del></del>		
<del></del> -		

OWNER Unisys Corporation, Great Neck, NY			
WELL NO. Recovery Well 1			
PAGE 1 OF 2 PAGES			
SCREEN TYPE Stainless-steel made by Cook Screen Company			
DIAM. 16-inch SLOT NO. 50			
SETTING 140-160 & 171-191/sump 191-196 ftbg			
SAND PACK Morie 2/3 mix			
CASING 16-inch black steel pipe			
SETTING 2.0.50 to 1/0.50 by			
SETTING 3.0 ft ag to 140 ft bg			
DEVELOPMENT			
DURATION			
STATIC WATER LEVEL			
YIELD			

REMARKS

DEPTH (FEET	л)   то	DESCRIPTION
0	30	SAND, medium to coarse: some fine to medium gravel: few cobbles and
		boulders: tan.
45	50	SILT and CLAY; some fine sand; gray.
	60	SILT and fine sand; gray.
	70	SAND, fine to very fine, some medium; little silt; gray.
	110	SAND, very fine to fine; tan.
	120	SAND, fine to very fine, trace medium; tan.
	130	SAND, fine to very fine, trace medium; trace silt and clay; tan.
	150	SAND, fine, some medium; tan; slightly coarser than 130 feet.
		SPLIT SPOONS
160	161.5	SAND, fine, some medium; some silt; orange, tan; 0.3-foot recovery.

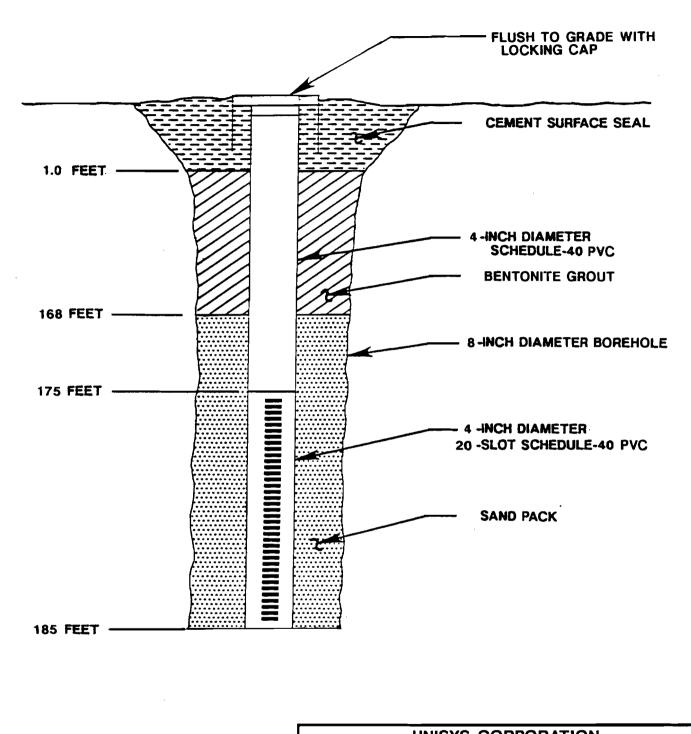
WELL NO	Recove	ery Well 1 PAGE 2 OF 2	PAGES
DEPTH (FEET)		DESCRIPTION	
180	181.5	SAND, fine to medium; little silt; tan-gray; 1.5-foot recovery	<del></del>
190	192	Driller notes smooth drilling, may be silt and clay.	
	193	Driller notes rough drilling may be coming out of silt and cla	y.
200	201.5	SAND, fine to medium; some silt; white; a little finer than 18	0 foot
<del></del>		sample; 1.3-foot recovery.	
210	213	Driller notes very smooth drilling (CLAY), lost circulation of	mud;
		white silty clay in mud.	
220	221.5	SAND, fine to very fine; some silt; orange-white; 0.7 foot thi	ck.
		SAND, fine to very fine with three 0.1 foot layers of white-or	ange
		silt and clay mixed in; 0.9 foot thick.	
<del> </del>	270	End of Borehole.	<del></del>
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GEOLOGIC LOG	OWNER Unisys Corporation, Great Neck, NY				
LEGGETTE, BRASHEARS & GRAHAM, INC.	WELL NO. Recovery Well 2				
WILTON, CONNECTICUT	PAGE 1 OF 2 PAGES				
LOCATION Reclamation area across	SCREEN TYPE Stainless-steel made by Cook Screen Company				
from Guard Post No. 3	DIAM. 8-inch SLOT NO. 30				
DATE COMPLETED July 19, 1991	<b>SETTING</b> screen 180-210/sump 210-215 ftbg				
DRILLING COMPANY R & L Well Drilling	SAND PACK Morie No. 2				
	CASING 8"I.D 8 3/4"O.D. black steel pipe				
DRILLING Mud rotary: 8" pilot hole METHOD reamed to 10 and 14"	CETTING 2 5 ft og to 190 ft ha				
SAMPLING 0 - 195 ft grab/wash samples	SETTING 2.5 ft ag to 180 ft bg				
METHOD 200 - 261 ft split-spoon samples	DEVELOPMENT				
OBSERVER John Benvegna	DURATION				
REFERENCE POINT (RP) Grade	STATIC WATER LEVEL				
ELEVATION OF RP	YIELD				
REMARKS					

DEPTH (FEET) FROM TO		DESCRIPTION			
	10	SAND, coarse to very coarse, and fine gravel; some cobbles.			
	28	SHEEN noticed on mud; odor at top of borehole; 6.5 ppm on PID.			
	35	SAND, coarse to very coarse; some fine gravel.			
	75	SAND, coarse to very coarse; some fine gravel; black sheen on mud.			
	95	SAND, coarse to medium, trace very coarse to fine; gray; slight odor.			
105	110	SILTY clay noted in strainer, bucket sample yielded medium to fine			
		sand.			
	115	Driller notes very dense material, bit rotation slowing; a lot of			
		brown silt and clay in mud.			
	115	SAND, fine to very fine; some silt; trace clay; gray.			
	130	Silty clay in mud; gray.			
	140	Driller notes medium sand in mud; coming out of silt and clay.			

OWNER	Unisys Corporation, Great Neck, New York					
WELL NO.	Recovery Well 2	PAGE	2	OF	2	PAGES

WELL NO.	Kecover	ry Well 2 PAGE 2 OF 2 PAGES
DEPTH (FE	ET)   TO	DESCRIPTION
	150	SAND, fine to very fine, little medium; trace silt; brown.
<del></del>	175	SAND, fine to very fine, some medium, little coarse; tan; gray.
<del></del>	185	SAND, medium to coarse, some very coarse; little silt and clay
		noted in mud.
<del></del>	195	SAND, fine to medium, little coarse, trace very fine; brown.
		SPLIT-SPOON SAMPLES
200	201.5	SAND, medium, some coarse, little very coarse; trace fine sand and
		gravel; brown; 0.4-foot recovery.
220	221.5	SAND, medium, some coarse; 0.3 foot layer of medium to coarse sand
		and fine gravel at bottom of spoon; 0.1 foot layer of red silt
		0.3 feet from bottom of spoon; 0.8-foot recovery.
240	241.5	SAND, coarse to medium, some very coarse; 1.10-foot recovery.
260	261.5	SAND, medium to fine, some very fine; little silt; streaks of black
		and brown silt throughout sample; 0.85-foot recovery.
		End of Borehole.
	<u> </u>	



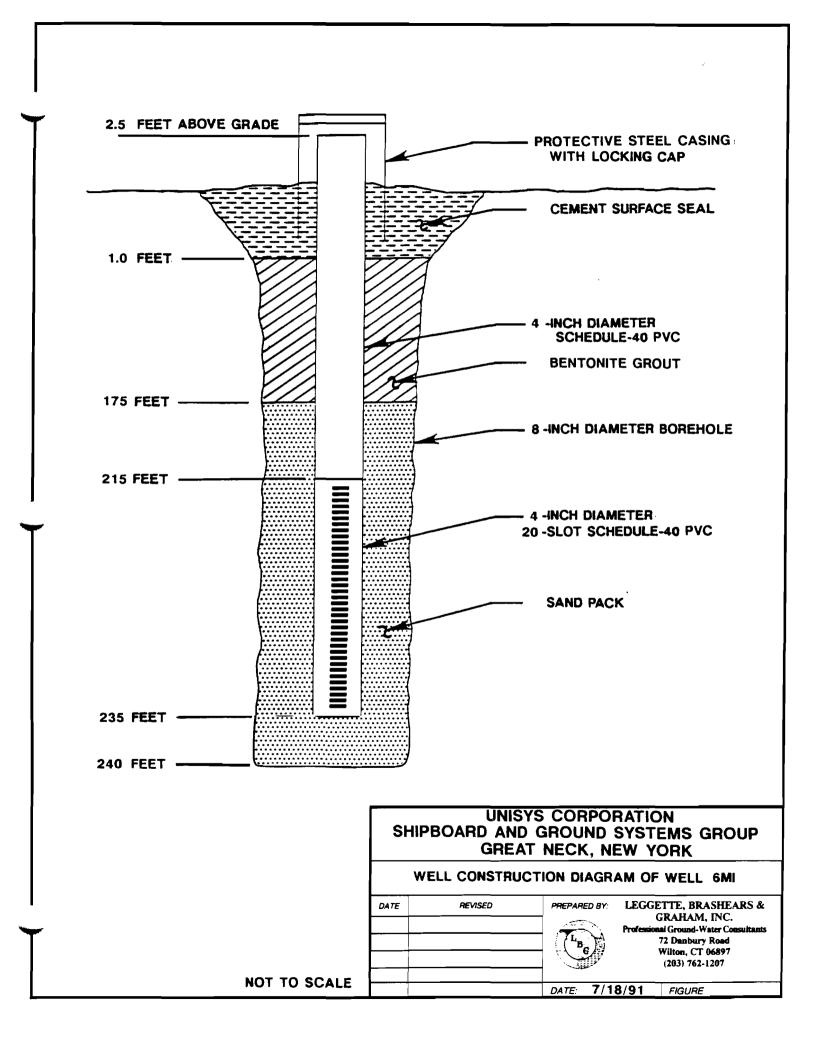
#### WELL CONSTRUCTION DIAGRAM OF WELL 2MU

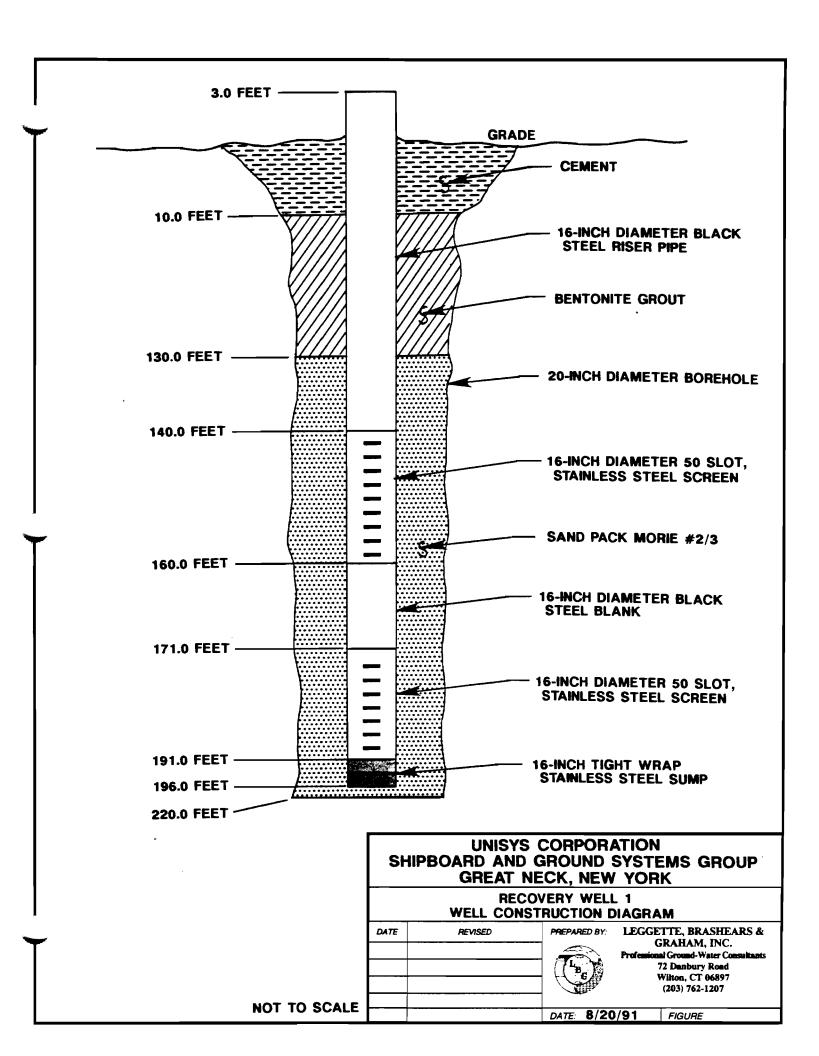
REVISED DATE PREPARED BY: 7/18/91 FIGURE DATE:

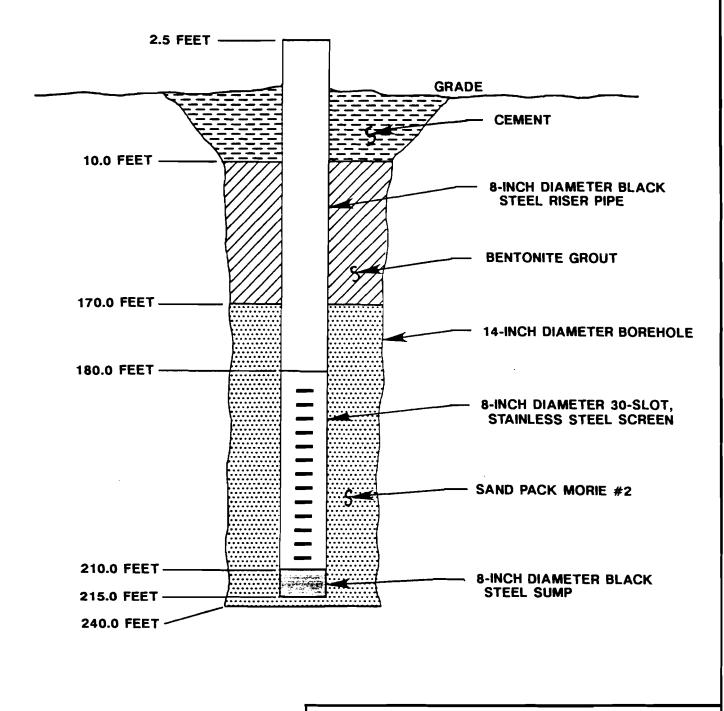
LEGGETTE, BRASHEARS & GRAHAM, INC. nal Ground-Water Consultants 72 Danbury Road Wilton, CT 06897

(203) 762-1207

**NOT TO SCALE** 







#### RECOVERY WELL 2 WELL CONSTRUCTION DIAGRAM

DATE REVISED PREPARED

PREPARED BY: LEGGETTE, BRASHEARS & GRAHAM, INC.
Professional Ground-Water Consultants
72 Danbury Road

rofessional Ground-Water Consultants 72 Danbury Road Wilton, CT 06897 (203) 762-1207

DATE: 8/20/91

FIGURE

NOT TO SCALE

# APPENDIX L5 LBG 1991 GROUND WATER CHEMISTRY



#### an environmental testing company

200 Monroe Turnpike Monroe, Connecticut 05468 (203) 261-4458 FAX (203) 268-5346

#### REPORT TRANSMITTAL

REPORT NUMBER 30910-1848

DATE October 22, 1991

CLIENT

Unisys Corporation 3199 Pilot Knob MS F1B05 Eagan, MN 55121

ATTENTION Mr. Kevin Krueger

The above referenced report is enclosed. Copies of this report and supporting data will be retained in our files in the event they are required for future reference.

If there are any questions concerning this report, please do not hesitate to contact us.

Any samples submitted to our Laboratory will be retained for a maximum of thirty (30) days from receipt of this report, unless other arrangements are desired.

30910-1848
UNISYS CORPORATION
3199 Pilot Knob
MS F1B05
Eagan, Minnesota 55121

Re: Great Neck, NY C-34141

Attention: Mr. Kevin Krueger

#### **PURPOSE**

Eighteen samples, four trip blanks and one field blank collected on September 17-20, 1991 were submitted to IEA, Inc. by Unisys Corporation. The client requested the samples be analyzed for TCL volatile organics plus a library search for non-target compounds.

#### METHODOLOGY

Volatile organics were determined using purge and trap GC/MS. The instrumentation used was a Tekmar Dynamic Headspace Concentrator interfaced with a Hewlett-Packard Model 5995C GC/MS/DS.

#### DISCUSSION

<u>Volatile Organics</u> - Sample VW-9 was run at a 1:100 dilution within holding time. This sample needed to be rerun at a 1:20 dilution which was done one day past holding time. Both sets of data have been reported. The 1:100 diluted run has been designated with the suffix "DL".

The following samples were diluted for target compounds:

<u>Sample</u>	<u>Dilution</u>	<u>Sample</u>	<u>Dilution</u>
6MI	1:5	VW-13	1:200
2MU	1:4	VW-8	1:100
8GL	1:2	VW-14	1:500
12MI	1:2	VW-14	1:500
10GL	1:2	VW-9*	1:20
7GL	1:2	VW-9 DL	1:100
5MI	1:2	2MI	1:5

<sup>\*</sup>out of holding time

The QC for sample VW-9 was run on the 8th day after receipt (one day past holding time). The client was contacted and requested this be noted in the case narrative and the data be reported.

The laboratory followed USEPA CLP-SOW Document #OLMO1.0 for the GC/MS calibration criteria.

#### **RESULTS**

The results are presented in the following Tables. Also enclosed is the data package containing all relevant QA/QC and raw data.

Prepared by:

Jeffrey C. Curran Laboratory Manager

JCC/adj

The liability of IEA, Inc. is limited to the actual dollar value of this project.

#### TABLE 1.0 30910-1848 UNISYS CORPORATION **EPA TCL VOLATILE ORGANICS**

All values are ug/L.

Sample Identification

<u>Dilution Factor</u>	1.0	1.0_	
Method_Blank I.D.	<u>&gt;A7848</u>	<u>&gt;A7848</u>	0 111 11
Compound	Method <u>Blank</u>	<u>8GU</u>	Quantitation Limits with no <u>Dilution</u>
Chloromethane Bromomethane Vinyl Chloride Chloroethane Methylene Chloride Acetone Carbon Disulfide 1,1-Dichloroethene 1,1-Dichloroethane 1,2-Dichloroethane 1,2-Dichloroethane 2-Butanone 1,1,1-Trichloroethane Carbon Tetrachloride Vinyl Acetate Bromodichloromethane 1,2-Dichloropropane cis-1,3-Dichloropropene Trichloroethene Dibromochloromethane 1,1,2-Trichloroethane Benzene trans-1,3-Dichloropropene Bromoform 4-Methyl-2-pentanone 2-Hexanone Tetrachloroethene 1,1,2,2-Tetrachloroethane Toluene Chlorobenzene Ethylbenzene	U U U U U U U U U U U U U U U U U U U	U U U U U U U U U U U U U U U U U U U	10 10 10 10 5 5 5 5 5 5 5 5 5 5 5 5 5 5
Styrene Xylene (total)	U U	U U	5 5

#### TABLE 1.1 30910-1848 UNISYS CORPORATION **EPA TCL\_VOLATILE ORGANICS**

All values are ug/L.

#### Sample Identification

<u>Dilution Factor</u>	1.0	1.0	1.0	<u>5.0</u>	100.0	100.0	
Method_Blank I.D.	<u>&gt;A7878</u>	>A7878	<u>&gt;A7878</u>	<u>&gt;A7878</u>	<u>&gt;A7878</u>	<u>&gt;A7878</u>	0 111 11
<u>Compound</u>	Method <u>Blank</u>	TB <u>09/21/91</u>	TB 09/24/91	<u>2MI</u>	<u>VW-8</u>	VW-9 DL	Quantitation Limits with no Dilution
Chloromethane	U	U	U	U	U	U	10
Bromomethane	U	U	U	U	U	U	10
Vinyl Chloride	U	U	U	U	U	U	10
Chloroethane	U	U	U	U	U	U	10
Methylene Chloride	3J	2JB	U	13JB	120JB	1 <b>70</b> JB	5
Acetone	41	<b>8</b> JB	5JB	<b>78</b> B	1,400B	3,800	10
Carbon Disulfide	U	U	U	IJ	U	U	5
1,1-Dichloroethene	U	U	U	U	U	U	5
1,1-Dichloroethane	U	U	U	U	U	U	5
1,2-Dichloroethene (total)	U	U	U	670	17,000	2,200	5
:hloroform	U	U	U	U	U	U	5
1,2-Dichloroethane	U	U	U	U	U	U	5
2-Butanone	U	U	U	U	U	U	10
1,1,1-Trichloroethane	U	U	U	U	U	U	5
Carbon Tetrachloride	U	U	U	U	U	U	5
Vinyl Acetate	U	U	U	U	U	U	10
Bromodichloromethane	U	U	U	U	U	U	5
1,2-Dichloropropane	U	U	U	U	U	U	5
cis-1,3-Dichloropropene	U	U	U	U	U	U	5
Trichloroethene	U	U	U	180	390J	620	5
Dibromochloromethane	U	U	U	U	U	U	5
1,1,2-Trichloroethane	U	U	U	U	U	U	5
Benzene	U	U	U	U	U	U	5
trans-1,3-Dichloropropene	U	U	U	U	U	U	5
Bromoform	U	U	U	U	U	U	5
4-Methy1-2-pentanone	U	U	U	U	U	U	10
2-Hexanone	U	U	IJ	U	U	U	10
Tetrachloroethene	U	U	U	100	830	680	5
1,1,2,2-Tetrachloroethane	U	U	U	U	U	U	5
Toluene	U	U	U	U	U	88J	5
Chlorobenzene	U	U	U	U	U	U	5
Ethylbenzene	U	U	U	U	U	U	5
Styrene	U	U	U	U	U	U	5
Xylene (total)	U	U	U	U	U	U	5

#### TABLE 1.2 30910-1848 UNISYS CORPORATION **EPA TCL VOLATILE ORGANICS**

All values are ug/L.

Sample Identification

<u>Dilution Factor</u>	1.0	1.0	_1.0_	2.0	2.0	1.0	
Method Blank I.D.	>A7858	_>A7858_	>A7858	>A7858	<u>&gt;A7858</u>	<u>&gt;A7858</u>	
Compound	Method <u>Blank</u>	TB <u>09/19/91</u>	TB <u>09/20/91</u>	_8GL_	_12MI_	_6GL	Quantitation Limits with no <u>Dilution</u>
Chloromethane	U	U	U	U	U	U	10
Bromomethane	U	U	U	U	U	U	10
Vinyl Chloride	Ü	Ü	Ū	U	U	U	10
Chloroethane	Ū	Ū	Ü	Ü	Ü	U	10
Methylene Chloride	Ü	1J	U	Ü	U	U	5
Acetone	<b>4</b> J	4JB	3JB	22B	6JB	6JB	10
Carbon Disulfide	U	U	U	U	U	U	5
1,1-Dichloroethene	U	U	Ü	U	2J	U	5
1,1-Dichloroethane	U	U	U	U	U	U	5
1,2-Dichloroethene (total)	ป	U	U	350	320	15	5
hloroform	U	2J	U	U	U	U	5
1,2-Dichloroethane	U	U	U	U	U	U	5
2-Butanone	U	U	U	U	U	U	10
1,1,1-Trichloroethane	U	U	U	U	Ų	U	5
Carbon Tetrachloride	U	U	U	U	U	U	5
Vinyl Acetate	U	U	U	U	U	U	10
Bromodichloromethane	U	U	U	U	U	U	5
1,2-Dichloropropane	U	U	U	U	U	U	5
cis-1,3-Dichloropropene	U	U	U	U	U	U	5
Trichloroethene	U	U	U	140	75	<b>3</b> J	5
Dibromochloromethane	U	U	U	U	U	U	5
1,1,2-Trichloroethane	U	U	U	U	U	U	5
Benzene	U	U	U	U	U	U	5
trans-1,3-Dichloropropene	U	U	U	U	U	U	5
Bromoform	U	U	U	U	U	U	5
4-Methyl-2-pentanone	U	U	U	U	U	U	10
2-Hexanone	U	U	U	U	IJ	U	10
Tetrachloroethene	U	U	U	94	73	2J	5
1,1,2,2-Tetrachloroethane	U	U	U	U	U	U	5
Toluene	U	U	U	U	U	U	5
Chlorobenzene	U	U	U	U	U	U	5
Ethylbenzene	U	U	U	U	U	U	5
Styrene	U	U	Ü	U	U	U	5
Xylene (total)	U	U	U	U	U	U	5

#### TABLE 1.3 30910-1848 UNISYS CORPORATION **EPA TCL VOLATILE ORGANICS**

All values are ug/L.

Sample Identification

<u>Dilution Factor</u>	1.0	1.0	_2.0_	1.0	2.0	2.0_	
Method Blank I.D.	>A7858	>A7858	>A7858	>A7858	<u>&gt;A7858</u>	<u>&gt;A7858</u>	
Compound	Method <u>Blank</u>	_12ML	10GL	_5GL	_7GL	_ <u>5MI</u>	Quantitation Limits with no <u>Dilution</u>
Chloromethane Bromomethane Vinyl Chloride Chloroethane	U U U	U U U	U U U	U U U	U U U	U U U	10 10 10 10
Methylene Chloride Acetone	Ŭ 4J	Ŭ 6JB	2J 14JB	Ŭ 1JB	Ŭ 13JB	Ŭ 11JB	5 10
Carbon Disulfide	U	U	U	U	U	Ü	5
1,1-Dichloroethene 1,1-Dichloroethane	U U	U U	U	U U	U U	U U	5
1,2-Dichloroethene (total) Chloroform	U U	41 U	11 <b>0</b> U	23 U	230 1J	300 U	5 5
1,2-Dichloroethane 2-Butanone	U U	U U	U U	U U	U U	U U	5 10
1,1,1-Trichloroethane Carbon Tetrachloride	U U	U U	2J U	U U	U U	U U	5 5
Vinyl Acetate Bromodichloromethane	U U	U U	U U	U U	U U	U U	10 5
<pre>1,2-Dichloropropane cis-1,3-Dichloropropene</pre>	U U	U U	U U	U U	U	U U	5 5
Trichloroethene Dibromochloromethane	U U	5 U	44 U	7 U	63 U	89 U	5 5
1,1,2-Trichloroethane Benzene	Ü	Ü	Ü	Ü U	Ü	Ü	5 5
trans-1,3-Dichloropropene Bromoform	Ü	Ŭ U	Ü	υ U	Ŭ V	Ü	5 5
4-Methyl-2-pentanone 2-Hexanone	Ŭ	Ü	Ŭ	Ü	Ü	Ŭ U	10 10
Tetrachloroethene 1,1,2,2-Tetrachloroethane	Ŭ	12 U	210 U	4J U	32 U	<b>4</b> 9 U	5 5
Toluene Chlorobenzene	บ U	Ü	Ü	Ü	ັ2J ປ	ŽJ U	5 5
Ethylbenzene Styrene	U U	U U	Ü	Ü	Ü	Ü	5 5
Xylene (total)	Ü	Ü	Ü	U	Ü	Ü	5

#### TABLE 1.4 30910-1848 UNISYS CORPORATION EPA TCL VOLATILE ORGANICS

All values are ug/L.

Sample Identification

<u>Dilution Factor</u>	1.0	_500.0_	1.0	200.0	20.0	
Method Blank I.D.	<u>&gt;A7932</u>	>A7932	<u>&gt;A7932</u>	<u>&gt;A7932</u>	<u>&gt;A7932</u>	
<u>Compound</u>	Method Blank		<u>VW-10</u>	<u>VW-13</u>	<u>VW-9</u>	Quantitation Limits with no <u>Dilution</u>
Chloromethane	U	U	U	U	IJ	10
Bromomethane	U	IJ	U	IJ	U	10
Vinyl Chloride	U	U	IJ	U	IJ	10
Chloroethane	U	IJ	U	IJ	U	10
Methylene Chloride	U	1,100J	IJ	U	Ŭ	5
Acetone	<b>6</b> J	16,000B	5JB	3,700B	380B	10
Carbon Disulfide	U	U	IJ	U	U	5
1,1-Dichloroethene	U	IJ	U	IJ	U	5
1,1-Dichloroethane	U	U	IJ	U	IJ	5
1,2-Dichloroethene (total)	U	IJ	30	18,000	2,100	5
hloroform	U	U	IJ	U	IJ	5
🛶 ,2-Dichloroethane	U	IJ	U	U	U	5
<b>2</b> -Butanone	U	U	IJ	U	ប	10
1,1,1-Trichloroethane	U	750J	1J	U	U	5
Carbon Tetrachloride	U	U	U	U	IJ	5
Vinyl Acetate	U	U	U	U	U	10
Bromodichloromethane	U	U	U	U	U	5
1,2-Dichloropropane	U	U	U	U	U	5
cis-1,3-Dichloropropene	U	U	U	U	U	5
Trichloroethene	U	98,000	64	220J	260	5
Dibromochloromethane	U	Ù	IJ	U	IJ	5
1,1,2-Trichloroethane	U	IJ	U	IJ	U	5
Benzene	U	U	IJ	U	U	5
trans-1,3-Dichloropropene	U	IJ	U	IJ	U	5
Bromoform	U	U	U	U	U	5
4-Methyl-2-pentanone	U	U	U	U	U	10
2-Hexanone	U	U	U	U	U	10
Tetrachloroethene	U	54,000	60	770J	430	5
<pre>1,1,2,2-Tetrachloroethane</pre>	U	U	U	U	U	5
Toluene	U	19,000	U	5,500	13J	5
Chlorobenzene	U	Ü	U	U	U	5
Ethylbenzene	U	<b>8</b> 50J	U	480J	U	5
Styrene	U	U	U	U	U	5
Xylene (total)	U	4,100	U	2,000	U	5

#### TABLE 1.5 30910-1848 UNISYS CORPORATION **EPA TCL VOLATILE ORGANICS**

All values are ug/L.

Sample Identification

<u>Dilution Factor</u>	1.0	20.0	20.0	
Method Blank I.D.	<u>&gt;B5675</u>	<u>&gt;B567<b>5</b></u>	<u>&gt;B5675</u>	
	Method	VW-9	VW-9	Quantitation Limits with no
Compound	<u>Blank</u>	<u>MS</u>	MSD	<u>Dilution</u>
Chloromethane	U	U	U	10
Bromomethane	U	U	U	10
Vinyl Chloride	U	U	U	10 10
Chloroethane	U	U	U U	10 5
Methylene Chloride Acetone	0 2J	U U	U	10
Carbon Disulfide	U	U	U	5
1,1-Dichloroethene	Ü	790X	840X	
1,1-Dichloroethane	Ü	U	U	5 5
1,2-Dichloroethene (total)	Ü	1,800X	2,000	5
hloroform	Ŭ	U	U U	5
1,2-Dichloroethane	Ü	Ŭ	Ŭ	5 5 5
2-Butanone	Ŭ	Ü	Ŭ	10
1,1,1-Trichloroethane	Ŭ	Ü	Ŭ	5
Carbon Tetrachloride	Ŭ	Ū	Ū	5
Vinyl Acetate	Ū	Ū	U	10
Bromodichloromethane	U	U	U	5
1,2-Dichloropropane	U	U	U	5
cis-1,3-Dichloropropene	U	U	U	5
Trichloroethene	U	1,200X	1,200X	5
Dibromochloromethane	U	U	U	5
1,1,2-Trichloroethane	U	U	U	5
Benzene	U	990X	1,000X	5 5 5 5
trans-1,3-Dichloropropene	U	U	U	
Bromoform	U	U	U	5
4-Methyl-2-pentanone	U	U	U	10
2-Hexanone	U	U	U 430	10 5
Tetrachloroethene 1,1,2,2-Tetrachloroethane	U U	390 U	430 U	5 5
Toluene	U	920X	980X	ى ت
Chlorobenzene	Ü	940X	960X	5
Ethylbenzene	Ü	U	U	5 5 5
Styrene	Ü	Ü	Ü	5
Xylene (total)	Ü	Ü	Ŭ	5
10 100041)	Ū	Ü	Ū	_

U, J, X - See Appendix for definition.

\_ote: Sample detection limit = quantitation limit x dilution factor.

#### TABLE 1.6 30910-1848 UNISYS CORPORATION EPA TCL VOLATILE ORGANICS

All values are ug/L.

Sample Identification

<u>Dilution Factor</u>	1.0	5.0_	1.0	4.0	_1.0	
Method Blank I.D.	<u>&gt;G6509</u>	>G6509	<u>&gt;G6509</u>	>G6509	<u>&gt;G6509</u>	
<u>Compound</u>	Method <u>Blank</u>	<u>6MI</u>	FB	<u>2MU</u>	TB 09/18/91	Quantitation Limits with no <u>Dilution</u>
Chloromethane	U	U	U	U	U	10
Bromomethane	U	U	U	U	U	10
Vinyl Chloride	U	U	U	U	U	10
Chloroethane	U	U	U	U	U	10
Methylene Chloride	U	5J	0.7J	5J	0.9J	5
Acetone	3J	67B	6JB	65B	4JB	10
Carbon Disulfide	U	Ü	U	U	U	5
1,1-Dichloroethene	Ü	Ü	Ü	U	U	5
1,1-Dichloroethane	Ū	Ũ	Ü	Ü	U	5
1,2-Dichloroethene (total)	Ŭ	600	Ŭ	670	Ü	5
Chloroform	Ũ	Ü	Ū	7J	<b>2</b> J	5
1,2-Dichloroethane	Ŭ	Ũ	Ũ	U	U	5
2-Butanone	Ū	Ū	Ü	Ū	Ü	10
1,1,1-Trichloroethane	Ü	Ü	Ŭ	Ŭ	Ü	5
Carbon Tetrachloride	Ū	Ū	Ŭ	Ū	Ü	5
Vinyl Acetate	ŭ	Ŭ	Ü	Ŭ	Ü	10
Bromodichloromethane	ŭ	Ü	Ü	Ü	Ü	5
1,2-Dichloropropane	Ü	Ŭ	Ŭ	Ŭ	Ü	5
cis-1,3-Dichloropropene	Ŭ	Ü	Ŭ	Ŭ	Ŭ	5
Trichloroethene	Ŭ	100	Ŭ	170	Ŭ	5
Dibromochloromethane	Ŭ	Ü	Ŭ	Ü	Ü	5
1,1,2-Trichloroethane	Ŭ	Ŭ	Ŭ	Ŭ	Ŭ	5
Benzene	Ŭ	Ŭ	Ŭ	Ŭ	Ŭ	5
trans-1,3-Dichloropropene	Ū	Ŭ	Ü	Ŭ	Ü	5
Bromoform	U	Ū	Ü	U	Ü	5
4-Methyl-2-pentanone	Ü	Ū	Ü	U	U	10
2-Hexanone	Ũ	Ū	Ū	Ū	Ü	10
Tetrachloroethene	Ü	110	Ū	120	Ü	5
1,1,2,2-Tetrachloroethane	Ū	Ü	Ü	Ü	Ü	5
Toluene	Ū	Ū	Ū	2J	U	5
Chlorobenzene	Ŭ	Ŭ	Ŭ	U	Ū	5
Ethylbenzene	Ŭ	Ū	Ü	Ü	Ū	5
Styrene	Ū	Ŭ	Ŭ	Ū	Ū	5
Xylene (total)	Ū	Ü	Ü	Ü	Ü	5

### TABLE 2.0 30910-1848 UNISYS CORPORATION VOLATILE TENTATIVELY IDENTIFIED COMPOUNDS

Sample Identification: Method Blank >A7848

CAS#	Compound RT	Estimated Concentration, ug/L
	None detected	
	Sample Identification: 8GU	
CAS#	CompoundRT	Estimated Concentration, ug/L
	None detected	
	Sample Identification: Method Blank	>A7878
CAS#	Compound RT	Estimated <u>Concentration, ug/L</u>
	None detected	
	Sample Identification: TB 09/21/	91
CAS#	Compound RT	Estimated Concentration, ug/L
	None detected	
	Sample Identification: TB 09/24/9	91
CAS#	Compound RT	Estimated Concentration, ug/L
	None detected	
	Sample Identification: 2MI	
CAS#	Compound RT	Estimated Concentration, ug/L
	None detected	

## TABLE 2.1 30910-1848 UNISYS CORPORATION VOLATILE TENTATIVELY IDENTIFIED COMPOUNDS

Sample Identification: VW-9 DL

CAS#	Compound	RT	Estimated Concentration, ug/L
	None detected		
	Sample Identification:	<b>VW-</b> 8	
CAS#	Compound	<u>RT</u>	Estimated Concentration, uq/L
	None detected		
	Sample Identification: Method	Blank >A	7858
CAS#	Compound	<u>RT</u>	Estimated Concentration, ug/L
	None detected		
	Sample Identification: TB	09/19/91	
CAS#	Compound	<u>RT</u>	Estimated Concentration, ug/L
	None detected		
	Sample Identification: TB	09/20/91	
CAS#	<u>Compound</u>	<u>RT</u>	Estimated Concentration, ug/L
	None detected		
	Sample Identification:	8GL	
CAS#	<u>Compound</u>	<u>RT</u>	Estimated Concentration, ug/L
	None detected		

## TABLE 2.2 30910-1848 UNISYS CORPORATION VOLATILE TENTATIVELY IDENTIFIED COMPOUNDS

Sample Identification: 12MI

CAS#	Compound	RT	Estimated Concentration, ug/L
	Unknown siloxane Unknown siloxane Unknown siloxane	22.23 25.29 22.42	35J 33J 14J
	Sample Identification:	6GL	
CAS#	Compound	<u>RT</u>	Estimated Concentration, ug/L
	Unknown siloxane Unknown alkane	22. <b>4</b> 1 26.28	44J 7J
	Sample Identification:	12ML	
CAS#	Compound	RT	Estimated Concentration, ug/L
	Unknown siloxane Unknown siloxane	25.30 22.43	21J 1 <b>4</b> J
	Sample Identification:	10GL	
CAS#	Compound	RT	Estimated Concentration, ug/L
	Unknown siloxane	22.36	29J
	Sample Identification:	5GL	
CAS#	Compound	RT	Estimated Concentration, ug/L
	Unknown siloxane	22.42	1 <b>1</b> J

 $<sup>\</sup>boldsymbol{J}$  - See Appendix for definition.

## TABLE 2.3 30910-1848 UNISYS CORPORATION VOLATILE TENTATIVELY IDENTIFIED COMPOUNDS

Sample Identification: 7GL

	,		
CAS#	Compound	RT	Estimated Concentration, ug/L
	None detected		
	Sample Identification:	5MI	
CAS#	Compound	RT	Estimated Concentration, ug/L
	None detected		
	Sample Identification: Method	Blank >A	7932
CAS#	Compound	<u>RT</u>	Estimated Concentration, ug/L
	None detected		
	Sample Identification:	VW-14	
CAS#	Compound	<u>RT</u>	Estimated Concentration, ug/L
	None detected		
	Sample Identification:	VW-10	
CAS#	Compound	RT	Estimated Concentration, ug/L
110543	Unknown siloxane Unknown siloxane Hexane	25.39 22.56 8.54	57J 7J 6J

J - See Appendix for definition.

### TABLE 2.4 30910-1848 UNISYS CORPORATION VOLATILE TENTATIVELY IDENTIFIED COMPOUNDS

Sample Identification: VW-13

<u>C</u> AS#		RT	Estimated Concentration, ug/L
	Unknown siloxane Unknown siloxane Unknown alkane Unknown siloxane	22.50 25.37 8.51 23.87	63,000J 48,00 <b>0</b> J 1,600J 1,000J
	Sample Identification:	<b>V</b> W-9	
CAS#	Compound	<u>RT</u>	Estimated Concentration, ug/L
1634-04-4	Methyl tert-butyl ether Unknown siloxane	8.05 25.39	450J 320J
110543	Hexane.	8.73	230J
	Sample Identification: Method	Blank >G	6509
CAS#	Compound	<u>RT</u>	Estimated Concentration, ug/L
	None detected		
	Sample Identification:	6MI	
CAS#	Compound	RT	Estimated Concentration, ug/L
76-13-1	Unknown alkane 1,1,2-Trichlorotrifluoroethane	8.66 6.44	70J 28J
70 13 1	1,1,2 Traciforder fraud dechane	0.44	200
	Sample Identification:	FB	
CAS#	Compound	RT	Estimated Concentration, ug/L
	Unknown alkane	8.70	5J

 $<sup>{\</sup>bf J}$  - See Appendix for definition.

### TABLE 2.5 30910-1848 UNISYS CORPORATION VOLATILE TENTATIVELY IDENTIFIED COMPOUNDS

Sample Identification: 2MU

CAS#	Compound	RT	Estimated Concentration, ug/L
76-13-1	Unknown alkane 1,1,2-Trichlorotrifluoroethane	8.67 6.29	100J 48J
	Sample Identification:	TB 09/18/91	
CAS#	Compound	RT	Estimated Concentration, ug/L
	Unknown alkane	8.64	6J

J - See Appendix for definition.

#### **APPENDIX**

- U Indicates that the compound was analyzed for but not detected.
- J Indicates that the compound was analyzed for and determined to be present in the sample. The mass spectrum of the compound meets the identification criteria of the method. The concentration listed is an estimated value, which is less than the specified minimum detection limit but is greater than zero.
- B This flag is used when the analyte is found in the blanks as well as the sample. It indicates possible sample contamination and warns the data user to use caution when applying the results of this analyte.
- N Indicates that the compound was analyzed for but not requested as an analyte. Value will not be listed on tabular result sheet.
- X Matrix spike compound.
- (1) Cannot be separated.
- (2) Decomposes to azobenzene. Measured and calibrated as azobenzene.
- A This flag indicates that a TIC is a suspected aldol condensation product.
- E Indicates that it exceeds calibration curve range.
- D This flag identifies all compounds identified in an analysis at a secondary dilution factor.
- C Confirmed by GC/MS.
- T Compound present in TCLP blank.



#### an environmental testing company -

200 Monroe Turnpike Monroe, Connecticut 06468 (203) 261-4458 FAX (203) 268-5346

#### REPORT TRANSMITTAL

REPORT NUMBER 30910-1848A

DATE October 31, 1991

CLIENT

Unisys Corporation 3199 Pilot Knob MS F1B05 Eagan, MN 55121

ATTENTION Mr. Kevin Krueger

The above referenced report is enclosed. Copies of this report and supporting data will be retained in our files in the event they are required for future reference.

If there are any questions concerning this report, please do not hesitate to contact us.

Any samples submitted to our Laboratory will be retained for a maximum of thirty (30) days from receipt of this report, unless other arrangements are desired.

30910-1848A UNISYS CORPORATION 3199 Pilot Knob MS F1B05 Eagan, Minnesota 55121

Re: Great Neck, NY

Attention: Mr. Kevin Krueger

#### **PURPOSE**

Twelve samples, two field blanks and three trip blanks were submitted to IEA, Inc. by Unisys Corporation. The client requested the samples be analyzed for TCL volatile organics plus a library search for non-target compounds.

#### **METHODOLOGY**

Volatile organics were determined using purge and trap GC/MS. The instrumentation used was a Tekmar Dynamic Headspace Concentrator interfaced with a Hewlett-Packard Model 5995C GC/MS/DS.

#### **DISCUSSION**

<u>Volatile Organics</u> - Samples 2GL and 11MI were analyzed undiluted within holding time. Analysis at the necessary dilutions (1:10 and 1:5 respectively) did not occur until after midnight of the last day of holding time. Both sets of data were reported. The matrix spike/matrix spike duplicate of sample 2GL were analyzed out of holding time.

The following samples were analyzed at dilutions due to high levels of target compounds:

<u>Sample</u>	<u>Dilution</u>
11GL	1:20
VW-12	1:10
VW-3	1:20
4MI	1:5
VW-2	1:1,000
4GL	1:2
3GL	1:5

#### **RESULTS**

The results are presented in the following Tables. Also enclosed are the data packages containing all relevant QA/QC and raw data.

Prepared by:

Jeffrey C. Curran

Laboratory Manager

JCC/mt

The liability of IEA, Inc. is limited to the actual dollar value of this project.

#### TABLE 1.0 30910-1848A UNISYS CORPORATION **EPA TCL VOLATILE ORGANICS**

All values are ug/L.

Sample Identification

<u>Dilution Factor</u>	1.0	_1.0	<u>1.0</u> _	1.0_	
Method Blank I.D.	<u>&gt;A7932</u>	_>A7932	<u>&gt;A7932</u>	<u>&gt;A7932</u>	Oventitetien
<u>Compound</u>	Method <u>Blank</u>	TB <u>09/25/91</u>	2GL MSB	_2GL	Quantitation Limits with no <u>Dilution</u>
Chloromethane Bromomethane Vinyl Chloride Chloroethane	U U U	U U U	U U U	U U U	10 10 10 10
Methylene Chloride Acetone Carbon Disulfide	บ 6J บ	บ บ ช	2J 20B U	U 11B U	5 10 5
1,1-Dichloroethene 1,1-Dichloroethane 2-Dichloroethene (total) Illianioroform	U U U	U U U	51X U U U	1J U 920E 1J	5 5 5 5
1,2-Dichloroethane 2-Butanone 1,1,1-Trichloroethane Carbon Tetrachloride	U U U	U U U	U U U	U U U	5 10 5 5
Vinyl Acetate Bromodichloromethane 1,2-Dichloropropane	Ü U U	U U U	U U U	U U U	10
cis-1,3-Dichloropropene Trichloroethene Dibromochloromethane 1,1,2-Trichloroethane	U U U	U U U	U 50X U U	U 170 U U	5 5 5 5 5
Benzene trans-1,3-Dichloropropene Bromoform 4-Methyl-2-pentanone	U U U U	Ü U U	54X U U U	U U U	5 5 5 10
2-Hexanone Tetrachloroethene 1,1,2,2-Tetrachloroethane	Ü U U	U U U	U U U	U 110 U	10 5 5
Toluene Chlorobenzene Ethylbenzene Styrene	U U U	U U U	46X 47X U U	U U U	5 5 5 5
Xylene (total)	Ŭ	Ü	Ü	Ü	5

#### TABLE 1.1 30910-1848A UNISYS CORPORATION **EPA TCL VOLATILE ORGANICS**

All values are ug/L.

Sample Identification

<u>Dilution Factor</u>	1.0	<u>10.0</u>	_1.0	1.0	
Method Blank I.D.	>A7932	>A7932	>A7932	>A7932	
<u>Compound</u>	Method <u>Blank</u>	2GL RE	FB <u>09/24/91</u>	_11MI_	Quantitation Limits with no <u>Dilution</u>
Chloromethane Bromomethane Vinyl Chloride Chloroethane Methylene Chloride Acetone Carbon Disulfide 1,1-Dichloroethene 1,1-Dichloroethene 1,1-Dichloroethene (total) Inloroform 1,2-Dichloroethane 2-Butanone 1,1,1-Trichloroethane Carbon Tetrachloride Vinyl Acetate Bromodichloromethane 1,2-Dichloropropane cis-1,3-Dichloropropene Trichloroethene Dibromochloromethane 1,1,2-Trichloroethane Benzene trans-1,3-Dichloropropene Bromoform 4-Methyl-2-pentanone 2-Hexanone Tetrachloroethene 1,1,2,2-Tetrachloroethane Toluene Chlorobenzene Ethylbenzene Styrene		UUUUU 250B UUUUUUUUUUUUUUUUUUUUUUUUUUUUUUUUUUUU		UUUUUUUUUUUUUUUUUUUUUUUUUUUUUUUUUUUUUU	10 10 10 10 5 10 5 5 5 5 5 5 5 5 5 5 5 5
Xylene (total)	Ŭ	Ü	Ü	Ü	5

#### TABLE 1.2 30910-1848A UNISYS CORPORATION **EPA TCL VOLATILE ORGANICS**

All values are ug/L.

Sample Identification

Compound   Blank   RE   11GL   9GL   Dilution	<u>Dilution Factor</u>	1.0	5.0	20.0	1.0	
Method   11MI	Method Blank I.D.	<u>&gt;A7932</u>	<u>&gt;A7932</u>	>A7932	<u>&gt;A7932</u>	0
Bromomethane	<u>Compound</u>			<u> 11GL</u>	9GL	Limits with no
2-Butanone       U       U       U       U       U       U       IJ       5         1,1,1-Trichloroethane       U       U       U       U       U       U       U       S         Vinyl Acetate       U       <	Bromomethane Vinyl Chloride Chloroethane Methylene Chloride Acetone Carbon Disulfide 1,1-Dichloroethene 1,1-Dichloroethane ,2-Dichloroethene (total) Chloroform	U U U U U U U U	U U U 110B U U 490	U U U 550B U U U 2,300 U	U U U 12B U 1J U 82 U	10 10 10 5 10 5 5 5
TrichÍoroethene       U       120       210       51       5         Dibromochloromethane       U       U       U       U       U       5         1,1,2-Trichloroethane       U       U       U       U       U       5         Benzene       U       U       U       U       U       U       5         trans-1,3-Dichloropropene       U       U       U       U       U       U       U       5         Bromoform       U       U       U       U       U       U       5         4-Methyl-2-pentanone       U       U       U       U       U       10         2-Hexanone       U       U       U       U       U       10         Tetrachloroethene       U       U       U       U       U       5         1,1,2,2-Tetrachloroethane       U       U       U       U       5         Chlorobenzene       U       U       U       U       5         Ethylbenzene       U       U       U       U       U       5         Styrene       U       U       U       U       U       U       U       U	2-Butanone 1,1,1-Trichloroethane Carbon Tetrachloride Vinyl Acetate Bromodichloromethane 1,2-Dichloropropane	U U U U U	U U U U U	U U U U U	U 1J U U U U	10 5 5 10
4-Methyl-2-pentanone       U       U       U       U       U       U       U       10         2-Hexanone       U       U       U       U       U       U       10         Tetrachloroethene       U       110       510       36       5       5         1,1,2,2-Tetrachloroethane       U       U       U       U       U       5         Toluene       U       4J       84J       U       5         Chlorobenzene       U       U       U       U       U         Ethylbenzene       U       U       U       U       U       5         Styrene       U       U       U       U       U       U       5	Trichloroethene Dibromochloromethane 1,1,2-Trichloroethane Benzene trans-1,3-Dichloropropene	U U U U	120 U U U U	210 U U U U	51 U U U U	5 5 5 5 5
Xylene (total) U U U 5	4-Methyl-2-pentanone 2-Hexanone Tetrachloroethene 1,1,2,2-Tetrachloroethane Toluene Chlorobenzene Ethylbenzene Styrene		U U 110 U 4J U U	U U 510 U 84J U U U	U 36 U U U U U	10 10 5 5 5 5 5

#### TABLE 1.3 30910-1848A UNISYS CORPORATION **EPA TCL VOLATILE ORGANICS**

All values are ug/L.

Sample Identification

<u>Dilution Factor</u>	1.0	10.0	10.0	2.0	<u>1.0</u>	
Method Blank I.D.	>B5675	>B5675	>B5675	<u>&gt;B5675</u>	_>B5675_	
Compound	Method <u>Blank</u>	2GL MS	2GL MSD	<u>4GL</u>	FB <u>09/25/91</u>	Quantitation Limits with no <u>Dilution</u>
Chloromethane	U	U	U	U	U	10
Bromomethane	U	U	U	U	U	10
Vinyl Chloride	U	U	U	U	U	10
Chloroethane	U	U	U	U	U	10
Methylene Chloride	U	U	U	U	U	5
Acetone	2J	210B	220B	6JB	5JB	10
Carbon Disulfide	U	U	U	U	U	5
1,1-Dichloroethene	U	430X	420X	U	U	5
1.1-Dichloroethane	U	U	U	U	U	5 5 5
2-Dichloroethene (total)	U	990	1,000	320	IJ	5
Chloroform	U	U	U	U	U	5
1,2-Dichloroethane	U	U	U	U	U	5
2-Butanone	U	U	U	U	3J	10
1,1,1-Trichloroethane	U	U	U	U	U	5
Carbon Tetrachloride	U	U	U	U	U	5
Vinyl Acetate	U	U	IJ	U	U	10
Bromodichloromethane	U	U	U	U	U	5
1,2-Dichloropropane	U	U	U	U	U	5
cis-1,3-Dichloropropene	U	U	U	IJ	U	5 5
Trichloroethene	U	590X	630X	91	U	5
Dibromochloromethane	U	U	U	U	U	5 5 5 5
1,1,2-Trichloroethane	U	U	U	U	U	5
Benzene	U	460X	490X	U	U	5
trans-1,3-Dichloropropene	U	U	U	U	U	5
Bromoform	U	IJ	U	ีป	U	5
4-Methyl-2-pentanone	U	U	U	U	U	10
2-Hexanone	U	U	U	U	U	10
Tetrachloroethene	U	120	120	58	U	5
1,1,2,2-Tetrachloroethane	U	U	U	U	U	5
Toluene	U	450X	460X	U	U	5
Chlorobenzene	บ	450X	470X	U	U	. 5 5 5
Ethylbenzene	U	U	U	U	U	5
Styrene	U	U	U	U	U	5
Xylene (total)	Ü	U	U	U	U	5

#### TABLE 1.4 30910-1848A UNISYS CORPORATION **EPA TCL VOLATILE ORGANICS**

All values are ug/L.

Sample Identification

<u>Dilution Factor</u>	1.0	<u>5.0</u>	10.0	1.0	1.0	
Method Blank I.D.	>B5683	>B5683	>B5683	>B5683_	>B5683	
<u>Compound</u>	Method <u>Blank</u>	<u>3GL</u>	<u><b>V</b>W-12</u>	TB <u>09/26/91</u>	TB <u>09/27/91</u>	Quantitation Limits with no <u>Dilution</u>
Chloromethane	U	U	U	U	U	10
Bromomethane	U	U	U	U	U	10
Vinyl Chloride	U	U	U	U	U	10
Chloroethane	U	U	U	U	U	10
Methylene Chloride	U	U	U	U	U	5
Acetone	<b>2</b> J	31JB	20JB	2JB	2JB	10
Carbon Disulfide	U	U	U	U	U	5
1,1-Dichloroethene	U	U	U	U	U	5
¹ 1-Dichloroethane	U	U	U	U	U	5
2-Dichloroethene (total)	U	820	1,200	U	U	5
chloroform	U	U	U	U	U	5
1,2-Dichloroethane	U	U	U	U	U	5
2-Butanone	U	U	U	U	U	10
1,1,1-Trichloroethane	U	U	U	U	U	5
Carbon Tetrachloride	U	U	U	U	U	5
Vinyl Acetate	U	U	U	U	U	10
Bromodichloromethane	U	U	U	U	U	5
1,2-Dichloropropane	U	U	U	U	U	5
cis-1,3-Dichloropropene	U	U	U	U	U ·	5
Trichloroethene	Ü	180	200	U	U	5
Dibromochloromethane	Ū	Ü	Ü	Ū	U	5
1,1,2-Trichloroethane	Ū	Ū	Ū	U	U	5
Benzene	Ü	U	U	U	U	5 5
trans-1,3-Dichloropropene	U	U	U	U	U	5
Bromoform	U	U	U	U	U	5
4-Methyl-2-pentanone	U	U	U	IJ	U	. 10
2-Hexanone	U	U	U	U	U	10
Tetrachloroethene	U	110	320	U	U	5
1,1,2,2-Tetrachloroethane	U	U	U	U	U	5
Toluene	U	U	U	U	5	5
Chlorobenzene	U	U	U	U	U	. 5
Ethylbenzene	U	Ū	Ū	U	Ü	5
Styrene	Ū	Ū	Ū	Ū	Ū	5 5
Xylene (total)	U	U	Ü	U	U	5

#### TABLE 1.5 30910-1848A UNISYS CORPORATION **EPA TCL VOLATILE ORGANICS**

All values are ug/L.

Sample Identification

<u>Dilution Factor</u>	1.0_	1.0	20.0	<u>5.0</u>	1,000.0	1.0	
Method Blank I.D.	>B5695	>B5695	>B5695	<u>&gt;B5695</u>	<u>&gt;B5695</u>	<u>&gt;B5695</u>	
Compound	Method <u>Blank</u>	<u>VW-11</u>	<u>VW-3</u>	<u>4MI</u>	<u>VW-2</u>	<u> VW-6</u>	Quantitation Limits with noDilution
Chloromethane	U	U	IJ	U	U	U	10
Bromomethane	ប	U	U	U	U	U	10
Vinyl Chloride	U	U	U	U	U	U	10
Chloroethane	U	U	U	U	บ	U	10
Methylene Chloride	U	U	U	U	U	U	5
Acetone	13	2JB	130JB	15JB	5,600JB	4JB	10
Carbon Disulfide	U	U	U	U	Ū	U	5
1,1-Dichloroethene	U	4J	U	U	U	U	5
1,1-Dichloroethane	U	U	U	U	U	U	5
2-Dichloroethene (total)	U	67	2,100	440	3,000J	44	5
Chloroform	U	U	Ú	U	໌ປ	U	5
1,2-Dichloroethane	U	U	U	U	U	U	5
2-Butanone	U	Ū	U	U	υ	U	10
1,1,1-Trichloroethane	U	4J	U	U	U	U	5
Carbon Tetrachloride	U	U	Ū	U	υ	U	5
Vinyl Acetate	Ū	Ū	Ū	Ũ	U	U	10
Bromodichloromethane	Ū	Ū	Ū	Ū	Ū	Ū	5
1,2-Dichloropropane	Ū	Ū	Ū	Ū	Ū	Ū	5
cis-1,3-Dichloropropene	บ	Ū	Ū	Ū	Ū	Ū	5
Trichloroethene	Ū	39	130	31	64,000	32	5
Dibromochloromethane	Ū	Ü	Ü	Ū	Ü	Ū	5
1,1,2-Trichloroethane	Ū	Ŭ	Ŭ	Ū	Ū	Ū	5
Benzene	Ū	Ū	U	U	U	U	5
trans-1,3-Dichloropropene	U	U	U	Ü	υ	U	5
Bromoform	U	Ü	U	U	บ	U	5
4-Methyl-2-pentanone	U	U	U	U	U	U	. 10
2-Hexanone	U	U	U	Ū	υ	U	10
Tetrachloroethene	U	30	310	U	21,000	21	5
1,1,2,2-Tetrachloroethane	U	IJ	U	U	Ú	U	5
Toluene	U	U	U	U	18,000	U	5
Chlorobenzene	Ū	Ŭ	Ū	Ū	Ü	Ū	5
Ethylbenzene	Ū	Ū	Ū	Ū	Ū	U	5
Styrene	Ū	Ū	Ū	U	Ū	U	5
Xylene (total)	Ü	Ü	Ü	Ü	2,300J	Ü	5

### TABLE 2.0 30910-1848A UNISYS CORPORATION VOLATILE TENTATIVELY IDENTIFIED COMPOUNDS

Sample Identification: Method Blank >A7932

CAS#		Compound	RT	Estimated Concentration, ug/L
	None detect	ed		
		Sample Identification:	TB 09/25/91	
CAS#		Compound	RT	Estimated Concentration, ug/L
110543	Hexane		8.54	6J
		Sample Identificati	on: 2GL	
CAS#		Compound	RT	Estimated Concentration, uq/L
110543	Hexane		8.57	25J
		Sample Identification	: 2GL RE	
CAS#		Compound	RT	Estimated Concentration, uq/L
110543	Hexane		8.54	110J
		Sample Identification:	FB 09/24/91	
CAS#		Compound	RT	Estimated Concentration, ug/L
	None detect	ed		

 ${\bf J}$  - See Appendix for definition.

## TABLE 2.1 30910-1848A UNISYS CORPORATION VOLATILE TENTATIVELY IDENTIFIED COMPOUNDS

Sample Identification: 11MI

CAS#	Compound	RT	Estimated <u>Concentration, ug/L</u>
110543	Unknown siloxane Hexane	22.79 8.77	47J 14J
	Sample Identification:	11MI RE	
CAS#	Compound	<u>RT</u>	Estimated Concentration, ug/L
110543 67641	Unknown siloxane Hexane 2-Propanone Unknown siloxane	22.51 8.59 6.25 25.37	53J 49J 27J 26J
	Sample Identification:	: 11GL	
CAS#	Compound	<u>RT</u>	Estimated Concentration, ug/L
110543	Hexane	8.59	16 <b>0</b> J
	Sample Identification	: 9GL	
CAS#	Compound	RT	Estimated Concentration, ug/L
110543 76131	Hexane 1,1,2-Trichlorotrifluoroethane	8.62 6.47	7J 9J
	Sample Identification: Metho	d Blank >B	5675
CAS#	Compound	RT	Estimated Concentration, uq/L
	None detected		

J - See Appendix for definition.

## TABLE 2.2 30910-1848A UNISYS CORPORATION VOLATILE TENTATIVELY IDENTIFIED COMPOUNDS

Sample Identification: 4GL

CAS#		Compound	_	<u>RT</u>	Estimated Concentration, ug/L
110543	Hexane			8.91	16J
		Sample Identification:	FB	09/25/91	
CAS#		Compound		RT	Estimated Concentration, uq/L
110543	Hexane			8.91	11J
		Sample Identificati	on:	3GL	
CAS#		Compound	_	RT	Estimated Concentration, uq/L
110543	Hexane			8.93	38J
		Sample Identificatio	n: '	<b>V</b> W-12	
CAS#		Compound		<u>RT</u>	Estimated Concentration, ug/L
110543	Hexane			8.91	1 <b>0</b> 0J
		Sample Identification:	ТВ	09/26/91	
CAS#		Compound		RT	Estimated Concentration, uq/L
110543	Hexane			8.95	11J

 $\boldsymbol{J}$  - See Appendix for definition.

## TABLE 2.3 30910-1848A UNISYS CORPORATION VOLATILE TENTATIVELY IDENTIFIED COMPOUNDS

Sample Identification: TB 09/27/91

CAS#	<u>Compound</u>	RT	Estimated Concentration, ug/L
110543	Hexane	8.93	12J
	Sample Identification: Method	Blank >B	5695
CAS#	<u>Compound</u>	RT	Estimated Concentration, ug/L
	None detected		
	Sample Identification:	VW-11	
CAS#	Compound	RT	Estimated Concentration, ug/L
	None detected		
	Sample Identification:	VW-3	
CAS#	Compound	RT	Estimated Concentration, ug/L
	None detected		
	Sample Identification:	4MI	
CAS#	Compound	<u>RT</u>	Estimated Concentration, ug/L
	None detected		. •

 ${\bf J}$  - See Appendix for definition.

### TABLE 2.4 30910-1848A UNISYS CORPORATION VOLATILE TENTATIVELY IDENTIFIED COMPOUNDS

Sample Identification: VW-2

CAS# Compound RT Concentration, uq/L

None detected

Sample Identification: VW-6

CAS# Compound RT Estimated
Concentration, uq/L

None detected

#### APPENDIX

- U Indicates that the compound was analyzed for but not detected.
- J Indicates that the compound was analyzed for and determined to be present in the sample. The mass spectrum of the compound meets the identification criteria of the method. The concentration listed is an estimated value, which is less than the specified minimum detection limit but is greater than zero.
- B This flag is used when the analyte is found in the blanks as well as the sample. It indicates possible sample contamination and warns the data user to use caution when applying the results of this analyte.
- N Indicates that the compound was analyzed for but not requested as an analyte. Value will not be listed on tabular result sheet.
- X Matrix spike compound.
- (1) Cannot be separated.
- (2) Decomposes to azobenzene. Measured and calibrated as azobenzene.
- A This flag indicates that a TIC is a suspected aldol condensation product.
- E Indicates that it exceeds calibration curve range.
- D This flag identifies all compounds identified in an analysis at a secondary dilution factor.
- C Confirmed by GC/MS.
- T Compound present in TCLP blank.



an environmental testing company -

200 Monroe Turnpike Monroe, Connecticut 06468 (203) 261-4458 FAX (203) 268-5346

#### REPORT TRANSMITTAL

DATE <u>October 31, 1991</u>

CLIENT

Unisys Corporation 3199 Pilot Knob MS F1B05 Eagan, MN 55121

ATTENTION Mr. Kevin Krueger

The above referenced report is enclosed. Copies of this report and supporting data will be retained in our files in the event they are required for future reference.

If there are any questions concerning this report, please do not hesitate to contact us.

Any samples submitted to our Laboratory will be retained for a maximum of thirty (30) days from receipt of this report, unless other arrangements are desired.

30910-1848B UNISYS CORPORATION 3199 Pilot Knobb MS F1B05 Eagan, Minnesota 55121

Re: Great Neck, NY C-34141

Attention: Mr. Kevin Krueger

#### **PURPOSE**

Ten samples, one field blank and three trip blanks collected on September 27, 30 and October 1, 1991 were submitted to IEA, Inc. by Unisys Corporation. The client requested the samples be analyzed for TCL volatile organics plus a library search for non-target compounds.

#### <u>METHODOLOGY</u>

Volatile organics were determined using purge and trap GC/MS. The instrumentation used was a Tekmar Dynamic Headspace Concentrator interfaced with a Hewlett-Packard Model  $5995C\ GC/MS/DS$ .

#### DISCUSSION

<u>Volatile Organics</u> - The continuing calibration on 10/05/91 met CLP 3/90 criteria, with the exception that chloroethane did not respond. This was caused by background interference of an ion similar to those ions present in chloroethane that made interpretation of the peak impossible. Samples FB, TB 09/28/91, 1MI, 13GL and 1GL were on last day of holding time so they were analyzed under this calibration.

The laboratory followed the USEPA CLP-SOW Document #OLMO1.0 for the GC/MS calibration criteria.

#### **RESULTS**

The results are presented in the following Tables.

Prepared by:

e**y &.**\ Curran axory\ Manager

JCC/adj

The liability of IEA, Inc. is limited to the actual dollar value of this project.

#### TABLE 1.0 30910-1848B UNISYS CORPORATION **EPA TCL VOLATILE ORGANICS**

All values are ug/L.

Sample Identification

<u>Dilution Factor</u>	1.0	_1.0	1.0	20.0	20.0	1.0	
Method Blank I.D.	>B5713	<u>&gt;B5713</u>	<u>&gt;B5713</u>	>B5713	<u>&gt;B5713</u>	<u>&gt;B5713</u>	0 - 1 1 1 1 1 1 1 1
<u>Compound</u>	Method Blank	<u>FB</u>	TB <u>09/28/91</u>	<u>1MI</u>	_13GL	1GL	Quantitation Limits with noDilution
Chloromethane	U	U	U	U	U	U	10
Bromomethane	U	U	U	U	U	U	10
Vinyl Chloride	U	U	U	U	U	U	10
Chloroethane	U	U	U	U	U	U	10
Methylene Chloride	U	U	U	U	U	U	5
Acetone	<b>4</b> J	5JB	U	590B	390B	U	10
Carbon Disulfide	U	U	U	U	U	Ü	5
1,1-Dichloroethene	U	U	U	U	U	U	5
1,1-Dichloroethane	U	U	U	U	U	U	5
1,2-Dichloroethene (total)	U	U	U	2,800	2,700	84	5
Chloroform	3ე	2JB	U	U	U	U	5
1,2-Dichloroethane	U	U	IJ	Ü	U	U	5
2-Butanone	U	U	U	U	U	U	10
1,1,1-Trichloroethane	U	U	U	U	U	U	5
Carbon Tetrachloride	U	U	U	U	U	U	5
Vinyl Acetate	U	U	U	U	U	U	10
Bromodichloromethane	U	U	U	U	U	U	5
1,2-Dichloropropane	U	U	U	U	U	U	5
cis-1,3-Dichloropropene	U	U	U	U	U	U	5
Trichloroethene	Ü	U	U	260	240	23	5
Dibromochloromethane	U	U	U	U	U	U	5
1,1,2-Trichloroethane	Ū	U	U	U	U	U	5
Benzene	Ŭ	Ü	Ü	Ü	U	U	5
trans-1,3-Dichloropropene	Ú	Ū	U	U	U	U	5
Bromoform	U	U	U	IJ	U	U	5
4-Methyl-2-pentanone	U	U	U	U	U	U	10
2-Hexanone	Ü	Ü	Ù	U	U	U	10
Tetrachloroethene	Ū	Ŭ	Ü	370	340	19	5
1,1,2,2-Tetrachloroethane	Ŭ	Ũ	Ū	U	Ü	U	5
Toluene	Ū	Ü	Ŭ	Ŭ	Ū	Ü	5
Chlorobenzene	Ŭ	Ü	Ŭ	Ū	Ü	Ū	5
Ethylbenzene	Ŭ	Ŭ	Ŭ	Ŭ	Ū	Ū	5
Styrene	Ŭ	Ū	Ū	Ũ	Ū	Ü	5
Xylene (total)	Ŭ	Ŭ	Ŭ	Ŭ	Ŭ	Ü	5
J ()	•	-					

#### TABLE 1.1 30910-1848B UNISYS CORPORATION EPA TCL VOLATILE ORGANICS

All values are ug/L.

Sample Identification

<u>Dilution Factor</u>	1.0	1.0	1.0	1.0	1.0	
Method_Blank I.D.	>B5723	>B5723	>B <b>5</b> 723	>B5723	<u>&gt;B5723</u>	
Compound	Method Blank		TB 10/01/91	1MI/L	1GU	Quantitation Limits with no <u>Dilution</u>
Chloromethane	U	U	U	U	U	10
Bromomethane	U	U	บ	U	U	10
Vinyl Chloride	U	U	U	U	U	10
Chloroethane	U	U	U	U	U	10
Methylene Chloride	U	5	U	U	U	5
Acetone	U	14	U	U	U	10
Carbon Disulfide	U	U	U	U	U	5
1,1-Dichloroethene	U	44X	U	U	U	5
1,1-Dichloroethane	U	U	U	U	U	5
1,2-Dichloroethene (total)	U	U	U	160	21	5 5
Chloroform	4J	U	2JB	Ü	U	5
بر_,2-Dichloroethane	U	U	U	U	U	5
2-Butanone	Ŭ	U	U	U	U	10
1,1,1-Trichloroethane	Ü	Ū	U	U	5	5
Carbon Tetrachloride	U	Ü	U	U	U	5
Vinyl Acetate	Ü	Ü	Ü	U	U	10
Bromodichloromethane	Ū	Ü	Ü	U	U	5
1,2-Dichloropropane	Ü	U	U	U	U	
cis-1,3-Dichloropropene	Ŭ	Ū	Ü	U	U	5 5
Trichloroethene	Ū	44X	Ŭ	190	4J	5
Dibromochloromethane	Ŭ	Ü	Ü	Ü	U	5
1,1,2-Trichloroethane	Ŭ	Ŭ	Ŭ	Ŭ	Ū	5
Benzene	Ū	48X	Ŭ	Ü	Ŭ	
trans-1,3-Dichloropropene	Ŭ	Ü	Ū	Ũ	U	5 5
Bromoform	Ü	Ū	Ü	Ŭ	Ū	5
4-Methyl-2-pentanone	Ū	Ū	Ū	U	U	10
2-Hexanone	Ū	Ü	Ū	Ú	U	10
Tetrachloroethene	Ü	Ū	Ū	37	U	5
1,1,2,2-Tetrachloroethane	Ū	Ŭ	Ü	Ü	U	5
Toluene	Ŭ	50X	Ü	Ü	Ŭ	5
Chlorobenzene	Ŭ	51X	Ü	Ū	Ū	
Ethylbenzene	Ü	Ü	ŭ	Ŭ	Ŭ	5 5
Styrene	Ŭ	Ŭ	Ŭ	Ü	Ŭ	5
Xylene (total)	ŭ	Ŭ	Ŭ	Ü	Ü	5 5

### TABLE 1.2 30910-1848B UNISYS CORPORATION EPA TCL VOLATILE ORGANICS

All values are ug/L.

Sample Identification

<u>Dilution Factor</u>	1.0	_1.0	<u>500.0</u>	1.0	
Method Blank I.D.	<u>&gt;B5723</u>	<u>&gt;B5723</u>	<u>&gt;B5723</u>	<u>&gt;B5723</u>	Quantitation
Compound	Method <u>Blank</u>		_ <b>VW</b> -5	1MI/L MS	Quantitation Limits with no <u>Dilution</u>
Chloromethane	U	U	U	U	10
Bromomethane	IJ	U	U	U	10
Vinyl Chloride	U	U	U	Ü	10
Chloroethane	IJ	U	U	U	10
Methylene Chloride	U	U	U	5	5
Acetone	U	U	1,100J	U	10
Carbon Disulfide	U	U	U	5	5
1,1-Dichloroethene	U	50X	U	55X	5
1,1-Dichloroethane	U	U	U	5	5
1,2-Dichloroethene (total)	U	160	48,000	160	5
Chloroform	<b>4</b> J	U	U	<b>6</b> B	5 5
,2-Dichloroethane	U	U	U	6	
2-Butanone	U	U	U	U	10
1,1,1-Trichloroethane	U	U	U	6	5
Carbon Tetrachloride	U	U	U	U	5
Vinyl Acetate	U	U	U	U	10
Bromodichloromethane	U	U	U	U	5
1,2-Dichloropropane	υ	U	U	5	5
cis-1,3-Dichloropropene	U	U	U	7	5
Trichloroethene	U	240XE	U	200X	5
Dibromochloromethane	U	U	U	3J	5
1,1,2-Trichloroethane	U	U	U	U	5 5 5
Benzene	U	52X	U	55X	5
trans-1,3-Dichloropropene	U	U	U	U	5
Bromoform	U	U	U	U	5
4-Methyl-2-pentanone	U	U	U	U	10
2-Hexanone	U	U	U	9J	10
Tetrachloroethene	U	36	U	32	5
1,1,2,2-Tetrachloroethane	U	U	U	U	5
Toluene	U	56X	8,700	59X	5
Chlorobenzene	U	<b>5</b> 5X	U	54X	5
Ethylbenzene	U	U	U	6	5
Styrene	U	U	U	5	5
Xylene (total)	U	U	U	15	5

#### TABLE 1.3 30910-1848B UNISYS CORPORATION **EPA TCL VOLATILE ORGANICS**

All values are ug/L.

Sample Identification

<u>Dilution Factor</u>	1.0	1.0_	1.0	500.0	500.0	
Method Blank I.D.	<u>&gt;B5744</u>	<u>&gt;B5744</u>	<u>&gt;B5744</u>	<u>&gt;B5744</u>	<u>&gt;B5744</u>	
<u>Compound</u>	Method <u>Blank</u>	<u>MW-13</u>	_1ML	<u>VW-4</u>	<u>VW-1</u>	Quantitation Limits with no <u>Dilution</u>
Chloromethane	U	U	U	U	U	10
Bromomethane	U	U	U	U	U	10
Vinyl Chloride	U	U	U	U	U	10
Chloroethane	U	U	U	U	U	10
Methylene Chloride	U	U	U	U	U	5
Acetone	U	U	U	9,100	U	10
Carbon Disulfide	U	U	U	U	U	5
1,1-Dichloroethene	U	U	U	U	U	5 5 5
1,1-Dichloroethane	U	U	IJ	U	U	5
1,2-Dichloroethene (total)	U	45	19	72,000	90,000	5
Chloroform	U	U	U	Ú	Ú	5
_1,2-Dichloroethane	U	U	U	U	U	5
2-Butanone	Ū	U	U	U	U	10
1,1,1-Trichloroethane	Ū	U	U	U	U	5
Carbon Tetrachloride	U	U	U	U	U	5
Vinyl Acetate	U	U	U	U	U	10
Bromodichloromethane	U	U	U	U	U	5
1,2-Dichloropropane	U	U	U	U	U	5
cis-1,3-Dichloropropene	U	U	U	U	U	5
Trichloroethene	U	10	10	U	U	5
Dibromochloromethane	Ū	Ū	Ü	U	U	5
1,1,2-Trichloroethane	Ū	Ū	U	U	U	5
Benzene	Ü	Ŭ	Ū	Ū	Ū	5
trans-1,3-Dichloropropene	Ū	Ū	Ū	Ū	U	5 5 5 5
Bromoform	Ū	Ū	Ū	U	U	5
4-Methyl-2-pentanone	Ū	Ū	Ū	Ū	U	10
2-Hexanone	Ū	Ū	Ü	U	U	10
Tetrachloroethene	Ū	12	17	U	11,000	5
1,1,2,2-Tetrachloroethane	U	IJ	U	U	Ú	5
Toluene	U	<b>3</b> J	<b>3</b> J	3,300	16,000	5
Chlorobenzene	Ū	U	U	Ú	Ú	5 5 5 5
Ethylbenzene	Ū	Ū	Ü	Ū	1,100J	5
Styrene	Ū	Ū	Ū	U	Ú	5
Xylene (total)	Ü	Ū	Ū	Ū	4,800	5

#### TABLE 1.4 30910-1848B UNISYS CORPORATION **EPA TCL VOLATILE ORGANICS**

All values are ug/L.

Sample Identification

<u>Dilution Factor</u>	1.0	1.0	
Method Blank I.D.	>A8082	_>A8082_	
<u>Compound</u>	Method <u>Blank</u>	TB <u>10/02/91</u>	Quantitation Limits with no <u>Dilution</u>
Compound  Chloromethane Bromomethane Vinyl Chloride Chloroethane Methylene Chloride Acetone Carbon Disulfide 1,1-Dichloroethene 1,2-Dichloroethene 1,2-Dichloroethene (total) Chloroform 1,2-Dichloroethane 2-Butanone 1,1,1-Trichloroethane Carbon Tetrachloride Vinyl Acetate Bromodichloromethane 1,2-Dichloropropane cis-1,3-Dichloropropene Trichloroethene Dibromochloromethane 1,1,2-Trichloroethane Benzene trans-1,3-Dichloropropene Bromoform 4-Methyl-2-pentanone 2-Hexanone Tetrachloroethene 1,1,2,2-Tetrachloroethane Toluene Chlorobenzene	Blank U U U U U I I U U U U U U U U U U U U	10/02/91  U U U U 13B U U U U U U U U U U U U U U U U U U U	10 10 10 10 5 10 5 5 5 5 5 5 5 5 5 5 5 5
Ethylbenzene Styrene	Ü U	Ü U	5 5 5 5
Xylene (total)	U	U	5

# TABLE 2.0 30910-1848B UNISYS CORPORATION VOLATILE TENTATIVELY IDENTIFIED COMPOUNDS

Sample Identification: Method Blank >B5713

CAS#	Compound RT	Estimated Concentration, ug/L
	None detected	
	Sample Identification: FB	
CAS#	CompoundRT	Estimated Concentration, ug/L
	None detected	
	Sample Identification: TB 09/28/91	
CAS#	CompoundRT	Estimated Concentration, ug/L
	None detected	
	Sample Identification: 1MI	
CAS#	Compound RT	Estimated Concentration, ug/L
	None detected	
	Sample Identification: 13GL	
CAS#	CompoundRT	Estimated Concentration, ug/L
	None detected	
	Sample Identification: 1GL	
CAS#	CompoundRT	Estimated Concentration, ug/L
	None detected	

## TABLE 2.1 30910-1848B UNISYS CORPORATION VOLATILE TENTATIVELY IDENTIFIED COMPOUNDS

Sample Identification: Method Blank >B5723

CAS#	Compound	RT	Estimated Concentration, ug/L
	None detected		
	Sample Identification: TB	10/01/91	
CAS#	Compound	<u>RT</u>	Estimated Concentration, uq/L
	None detected		
	Sample Identification:	1MI/L	
CAS#	Compound	RT	Estimated Concentration, ug/L
<b>7</b> 6131	1,1,2-Trichlorotrifluoroethane	6.90	5J
	Sample Identification:	1GU	
CAS#	Compound	<u>RT</u>	Estimated Concentration, ug/L
	None detected		
	Sample Identification:	VW-5	
CAS#	Compound	RT	Estimated Concentration, uq/L
	None detected		
	Sample Identification: Method E	81ank >B5	744
CAS#	Compound	<u>RT</u>	Estimated Concentration, ug/L
	None detected		

J - See Appendix for definition.

## TABLE 2.2 30910-1848B UNISYS CORPORATION VOLATILE TENTATIVELY IDENTIFIED COMPOUNDS

Sample Identification: MW-13

CAS#	Compound None detected .	<u>RT</u>	Estimated Concentration, ug/L
	Sample Identification:	1ML	
CAS#	Compound	<u> RT</u>	Estimated Concentration, ug/L
	None detected		
	Sample Identification:	VW-4	
CAS#	Compound	RT	Estimated Concentration, ug/L
	None detected		
	Sample Identification:	VW-1	
CAS#	Compound	<u>RT</u>	Estimated Concentration, ug/L
	None detected		
	Sample Identification: Method	Blank >A8	3082
CAS#	Compound	RT	Estimated Concentration, ug/L
	None detected		
	Sample Identification: TB	10/02/91	
CAS#	<u>Compound</u>	<u>RT</u>	Estimated Concentration, ug/L
	None detected		

#### APPENDIX

- U Indicates that the compound was analyzed for but not detected.
- J Indicates that the compound was analyzed for and determined to be present in the sample. The mass spectrum of the compound meets the identification criteria of the method. The concentration listed is an estimated value, which is less than the specified minimum detection limit but is greater than zero.
- B This flag is used when the analyte is found in the blanks as well as the sample. It indicates possible sample contamination and warns the data user to use caution when applying the results of this analyte.
- N Indicates that the compound was analyzed for but not requested as an analyte. Value will not be listed on tabular result sheet.
- X Matrix spike compound.
- (1) Cannot be separated.
- (2) Decomposes to azobenzene. Measured and calibrated as azobenzene.
- A This flag indicates that a TIC is a suspected aldol condensation product.
- E Indicates that it exceeds calibration curve range.
- D This flag identifies all compounds identified in an analysis at a secondary dilution factor.
- C Confirmed by GC/MS.
- T Compound present in TCLP blank.

### APPENDIX K

### ROUX ASSOCIATES AQUIFER PERFORMANCE TEST