

**JUNE 2001 QUARTERLY
GROUNDWATER MONITORING REPORT
FOR THE
WEISENBERGER TIE & LUMBER SITE
MARATHON CITY, WISCONSIN**

October 30, 2001

Robert E. Lee & Associates, Inc.

Engineering • Surveying • Laboratory Service
2825 South Webster Avenue

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Robert E. Lee & Associates, Inc.
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October 30, 2001

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Mr. John Grump
WISCONSIN DEPARTMENT OF NATURAL RESOURCES
P.O. Box 4001
Eau Claire, WI 54702

RE: June 2001 Quarterly Groundwater Monitoring Report
Weisenberger Tie and Lumber Company
WDNR File #95S440

Dear Mr. Grump:

Please find enclosed the quarterly groundwater monitoring report for the above-named site for the sampling event of June 27, 2001. The following items require comment:

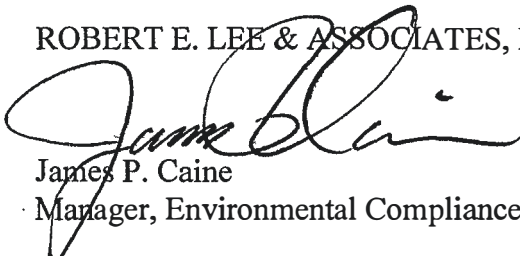
- A sample could not be collected from monitoring well DMW-3, since it was dry.

The results of the groundwater sampling continue to identify several enforcement standard exceedances; however, the groundwater plume appears to be relatively stable.

If you have any questions and/or comments regarding this matter, please contact our office.

Sincerely,

ROBERT E. LEE & ASSOCIATES, INC.



James P. Caine
Manager, Environmental Compliance
JPC/njm

ENC.

**MONITORING WELLS
DIOXIN/FURAN ANALYSIS**

WEISENBERGER TIE & LUMBER COMPANY

Dioxin/Furan Analytical Results - MW-2

Date Sampled	Compound (pg/l)																I-TEQ/89 2,3,7,8-TCDD	
	2,3,7,8-TCDD	1,2,3,7,8-PeCDD	1,2,3,4,7,8-HxCDD	1,2,3,6,7,8-HxCDD	1,2,3,7,8,9-HxCDD	1,2,3,4,6,7,8-HpCDD	1,2,3,4,6,7,8,9-OCDD	2,3,7,8-TCDF	1,2,3,7,8-PeCDF	2,3,4,7,8-PeCDF	1,2,3,4,7,8-HxCDF	1,2,3,6,7,8-HxCDF	2,3,4,6,7,8-HxCDF	1,2,3,7,8,9-HxCDF	1,2,3,4,6,7,8-HpCDF	1,2,3,4,7,8,9-HpCDF		1,2,3,4,6,7,8,9-OCDF
I-TEF/89	1.00	0.50	0.10	0.10	0.10	0.01	0.001	0.10	0.05	0.50	0.10	0.1	0.1	0.1	0.01	0.01	0.001	3.00
1/19/1993	0	0	0	0	0	31	489	0	0	0	0	0	3.1	0	0	0	0	1.11

I-TEF/89 = International Toxicity Equivalent Factors/1989

I-TEQ/89 = International Toxicity Equivalents (based on I-TEF/89)

= Exceedance of I-TEQ/89

= Results are above detection limit but below quantitation limit

WEISENBERGER TIE & LUMBER COMPANY

Dioxin/Furan Analytical Results - MW-3

Date Sampled	Compound (pg/l)																I-TEQ/89 2,3,7,8-TCDD	
	2,3,7,8-TCDD	1,2,3,7,8-PeCDD	1,2,3,4,7,8-HxCDD	1,2,3,6,7,8-HxCDD	1,2,3,7,8,9-HxCDD	1,2,3,4,6,7,8-HpCDD	1,2,3,4,6,7,8,9-OCDD	2,3,7,8-TCDF	1,2,3,7,8-PeCDF	2,3,4,7,8-PeCDF	1,2,3,4,7,8-HxCDF	1,2,3,6,7,8-HxCDF	2,3,4,6,7,8-HxCDF	1,2,3,7,8,9-HxCDF	1,2,3,4,6,7,8-HpCDF	1,2,3,4,7,8,9-HpCDF		1,2,3,4,6,7,8,9-OCDF
I-TEF/89	1.00	0.50	0.10	0.10	0.10	0.01	0.001	0.10	0.05	0.50	0.10	0.1	0.1	0.1	0.01	0.01	0.001	3.00
8/20/1992	0	0	12.4	528	44	11970	71770	8.50	24	20	66	24.4	43.8	0	970	82	6190	291.91
12/23/1999	<1.5	<2.2	<3.6	19	<3.3	410	3500	<1.1	<2.3	<1.6	<2.8	<3.5	3.7	<4.4	<3.2	<4.6	300	10.17
3/6/2000	<8.0	<8.0	<12	28	<9.6	500	4200	10	<9.4	<6.6	<9.3	<9.2	<13	<8.2	41	<9.4	370	13.78
6/30/2000	<3.2	<9.2	<19	37	<22	590	5200	<2.4	<3.0	<8.8	<13	<8.3	<5.7	<7.2	69	<11	420	15.91
9/27/2000	<8.1	<7.3	<11	13	<11	190	1800	<3.9	<4.9	<5.0	<4.1	11	<3.8	<5.0	<8.9	<10	94	6.19
3/28/2001	<4.2	<4.1	5.3	36	<1.9	710	5700	<3.1	<2.7	<2.7	9.2	<3.0	<1.6	<1.4	69	5.9	510	19.11
06/27/2001	<4.7	<9.5	<9.5	12.0	<9.5	220	2000	<3.0	<9.5	<9.5	<9.5	<9.5	<9.5	<9.5	14	<14	170	5.71

I-TEF/89 = International Toxicity Equivalent Factors/1989

I-TEQ/89 = International Toxicity Equivalents (based on I-TEF/89)

= Exceedance of I-TEQ/89

= Results are above detection limit but below quantitation limit

WEISENBERGER TIE & LUMBER COMPANY

Dioxin/Furan Analytical Results - MW-5

Date Sampled	Compound (pg/l)																I-TEQ/89 2,3,7,8-TCDD	
	2,3,7,8-TCDD	1,2,3,7,8-PeCDD	1,2,3,4,7,8-HxCDD	1,2,3,6,7,8-HxCDD	1,2,3,7,8,9-HxCDD	1,2,3,4,6,7,8-HpCDD	1,2,3,4,6,7,8,9-OCDD	2,3,7,8-TCDF	1,2,3,7,8-PeCDF	2,3,4,7,8-PeCDF	1,2,3,4,7,8-HxCDF	1,2,3,6,7,8-HxCDF	2,3,4,6,7,8-HxCDF	1,2,3,7,8,9-HxCDF	1,2,3,4,6,7,8-HpCDF	1,2,3,4,7,8,9-HpCDF		1,2,3,4,6,7,8,9-OCDF
I-TEF/89	1.00	0.50	0.10	0.10	0.10	0.01	0.001	0.10	0.05	0.50	0.10	0.1	0.1	0.1	0.01	0.01	0.001	3.00
8/20/1992	0	0	0	1.3	0	34.6	685	0	0	0	0	0	3.0	0	0	0	5.5	1.47
6/30/2000	<4.3	<8.5	<5.6	<5.6	<5.1	16	240	<4.1	<3.7	<3.9	<3.8	<3.6	<3.4	<4.4	<4.1	<5.5	<9.7	0.40
9/27/2000	<10	<5.2	<3.8	<5.7	<3.5	12	150	<5.6	<3.6	<2.9	<3.2	<2.4	<4.1	<4.6	<4.1	<3.4	<8.9	0.27
12/27/2000	<9.1	<2.8	<2.1	<3.2	<2.6	5.2	250	<5.8	<3.6	<2.6	<2.1	<2.3	<2.4	<3.0	6.5	<4.1	9.6	0.38
03/28/2001	<4.7	<2.6	<2.8	<3.1	<2.0	8.4	300	<3.6	<2.5	<1.7	<1.8	<2.4	<2.3	<1.6	<1.2	<1.2	<2.3	0.38
06/27/2001	<6.1	<9.5	<9.5	<9.5	<9.5	<16	130	<4.6	<9.5	<9.5	<9.5	<9.5	<9.5	<9.5	<13	<11	<39	0.13

I-TEF/89 = International Toxicity Equivalent Factors/1989

I-TEQ/89 = International Toxicity Equivalent (based on I-TEF/89)

= Exceedance of I-TEQ/89

= Results are above detection limit but below quantitation limit

WEISENBERGER TIE & LUMBER COMPANY

Dioxin/Furan Analytical Results - MW-6

Date Sampled	Compound (pg/l)																I-TEQ/89 2,3,7,8-TCDD	
	2,3,7,8-TCDD	1,2,3,7,8-PeCDD	1,2,3,4,7,8-HxCDD	1,2,3,6,7,8-HxCDD	1,2,3,7,8,9-HxCDD	1,2,3,4,6,7,8-HpCDD	1,2,3,4,6,7,8,9-OCDD	2,3,7,8-TCDF	1,2,3,7,8-PeCDF	2,3,4,7,8-PeCDF	1,2,3,4,7,8-HxCDF	1,2,3,6,7,8-HxCDF	2,3,4,6,7,8-HxCDF	1,2,3,7,8,9-HxCDF	1,2,3,4,6,7,8-HpCDF	1,2,3,4,7,8,9-HpCDF		1,2,3,4,6,7,8,9-OCDF
I-TEF/89	1.00	0.50	0.10	0.10	0.10	0.01	0.001	0.10	0.05	0.50	0.10	0.1	0.1	0.1	0.01	0.01	0.001	3.00
1/19/1993	0	0	0	0	0	39	763	0	0	0	0	0	3.9	0	0	0	0	1.55
6/30/2000	<0.75	<0.41	<0.62	7.2	2.7	100	1100	<0.6	<0.46	<0.59	<0.64	<0.41	1.3	<0.24	19	<2.0	110	3.52
9/27/2000	<9.4	<9.9	<13	<14	<6.8	120	1200	<7.9	<7.0	<4.2	<13	18	<11	<4.6	17	<8.0	120	4.49
12/27/2000	<4.9	<2.5	<3.0	51	6.5	1200	12000	<3.3	4.8	<1.4	14.0	<4.1	<1.2	<1.4	<2.1	17	1500	33.06
03/28/2001	<6.1	<2.6	19	95	9.6	2100	21000	<4.0	<9.0	<2.7	41	12	10	7.1	290	26	2300	66.83
06/27/2001	<8.1	<9.5	<9.5	<9.5	<9.5	200	2100	<5.2	<9.5	<9.5	<9.5	<9.5	<9.5	<9.5	450	<16	240	8.84

I-TEF/89 = International Toxicity Equivalent Factors/1989

I-TEQ/89 = International Toxicity Equivalents (based on I-TEF/89)

= Exceedance of I-TEQ/89

= Results are above detection limit but below quantitation limit

WEISENBERGER TIE & LUMBER COMPANY

Dioxin/Furan Analytical Results - MW-7

Date Sampled	Compound (pg/l)																I-TEQ/89 2,3,7,8-TCDD	
	2,3,7,8-TCDD	1,2,3,7,8-PeCDD	1,2,3,4,7,8-HxCDD	1,2,3,6,7,8-HxCDD	1,2,3,7,8,9-HxCDD	1,2,3,4,6,7,8-HpCDD	1,2,3,4,6,7,8,9-OCDD	2,3,7,8-TCDF	1,2,3,7,8-PeCDF	2,3,4,7,8-PeCDF	1,2,3,4,7,8-HxCDF	1,2,3,6,7,8-HxCDF	2,3,4,6,7,8-HxCDF	1,2,3,7,8,9-HxCDF	1,2,3,4,6,7,8-HpCDF	1,2,3,4,7,8,9-HpCDF		
I-TEF/89	1.00	0.50	0.10	0.10	0.10	0.01	0.001	0.10	0.05	0.50	0.10	0.1	0.1	0.1	0.01	0.01	0.001	3.00
12/23/1999	<3.0	<2.5	<2.3	12.0	<2.2	230	2100	<1.7	<1.0	<1.5	<2.0	<2.0	4.50	<1.9	25	<2.9	240	6.54
3/6/2000	<3.9	<5.4	<14	230	16	530	70000	<2.8	<11	<4.2	<14	<9.4	<10	<4.9	610	56	7200	113.76
6/30/2000	<0.34	4.9	3.5	20.0	5.9	400	5600	<0.43	1.1	<0.35	<0.41	1.5	3.4	2.0	55	7.6	480	16.84
9/27/2000	NO SAMPLE																	
12/27/2000	<3.9	<2.4	<2.0	60.0	4.5	1500	15000	<2.0	3.0	5.3	10.0	<3.8	9.3	5.2	<0.86	14	1300	43.14
03/28/2001	NO SAMPLE																	
06/27/2001	<6.4	<9.5	<9.5	21.0	<9.5	610	6600	<2.9	<9.5	<9.5	<9.5	<10	<9.5	<9.5	720	<14	510	22.51

I-TEF/89 = International Toxicity Equivalent Factors/1989

I-TEQ/89 = International Toxicity Equivalents (based on I-TEF/89)

 = Exceedance of I-TEQ/89

 = Results are above detection limit but below quantitation limit

WEISENBERGER TIE & LUMBER COMPANY

Dioxin/Furan Analytical Results - MW-10

Date Sampled	Compound (pg/l)																	I-TEQ/89 2,3,7,8-TCDD
	2,3,7,8-TCDD	1,2,3,7,8-PeCDD	1,2,3,4,7,8-HxCDD	1,2,3,6,7,8-HxCDD	1,2,3,7,8,9-HxCDD	1,2,3,4,6,7,8-HpCDD	1,2,3,4,6,7,8,9-OCDD	2,3,7,8-TCDF	1,2,3,7,8-PeCDF	2,3,4,7,8-PeCDF	1,2,3,4,7,8-HxCDF	1,2,3,6,7,8-HxCDF	2,3,4,6,7,8-HxCDF	1,2,3,7,8,9-HxCDF	1,2,3,4,6,7,8-HpCDF	1,2,3,4,7,8,9-HpCDF	1,2,3,4,6,7,8,9-OCDF	
I-TEF/89	1.00	0.50	0.10	0.10	0.10	0.01	0.001	0.10	0.05	0.50	0.10	0.1	0.1	0.1	0.01	0.01	0.001	3.00
1/19/1993	0	0	244	183	12.3	3190	18380	3.50	8.1	6.2	23.5	7.0	17.4	3.3	148	9.20	541	105.30
12/23/1999	<3.5	<6.4	110	6100	240	82000	500000	<9.9	290	210	510	<8.4	220	330	5600	300	21000	2270.50
3/6/2000	<7.5	<7.3	<25	2200	120	35000	240000	<11	<8.0	99	<33	<35	120	23	2100	160	10000	918.40
6/30/2000	1.1	7.9	11	170	17	2100	16000	2.2	7.1	12	<0.73	7.0	6.8	10	160	10	570	73.08
9/27/2000	<2.7	<3.1	<3.4	47	<2.0	700	4500	<1.8	<1.6	<2.0	4.3	4.8	<5.1	<1.2	44	<1.6	130	17.68
12/27/2000	<2.2	6.2	12	2700	110	34000	190000	38	92	170	280	<0.47	200	170	<1.1	110	4000	978.80
03/28/2001	NO SAMPLE																	
06/27/2001	<5.3	<9.5	<9.5	260	25	4200	26000	5.6	15	21	35	<9.5	21	22	730	<14	380	123.79

I-TEF/89 = International Toxicity Equivalent Factors/1989

I-TEQ/89 = International Toxicity Equivalents (based on I-TEF/89)

 = Exceedance of I-TEQ/89

 = Results are above detection limit but below quantitation limit

WEISENBERGER TIE & LUMBER COMPANY

Dioxin/Furan Analytical Results - DMW-1

Date Sampled	Compound (pg/l)																I-TEQ/89 2,3,7,8-TCDD	
	2,3,7,8-TCDD	1,2,3,7,8-PeCDD	1,2,3,4,7,8-HxCDD	1,2,3,6,7,8-HxCDD	1,2,3,7,8,9-HxCDD	1,2,3,4,6,7,8-HpCDD	1,2,3,4,6,7,8,9-OCDD	2,3,7,8-TCDF	1,2,3,7,8-PeCDF	2,3,4,7,8-PeCDF	1,2,3,4,7,8-HxCDF	1,2,3,6,7,8-HxCDF	2,3,4,6,7,8-HxCDF	1,2,3,7,8,9-HxCDF	1,2,3,4,6,7,8-HpCDF	1,2,3,4,7,8,9-HpCDF		1,2,3,4,6,7,8,9-OCDF
I-TEF/89	1.00	0.50	0.10	0.10	0.10	0.01	0.001	0.10	0.05	0.50	0.10	0.1	0.1	0.1	0.01	0.01	0.001	3.00
8/20/1992	0	0	0	1010	103	33320	248820	14.3	66.1	39.3	181	63.1	212	187	2610	416	21540	834
3/6/2000	<6.4	37	97	5200	290	100000	850000	<5.0	<3.1	<3.4	<14	<17	350	73	11000	1000	90000	2680
6/30/2000	<9.4	<5.5	510	19000	970	190000	150000	140	<48	980	1700	590	<19	1400	41000	3800	60000	5479
9/27/2000	<8.8	<7.7	<7.5	3300	140	71000	610000	32	<1.6	250	590	150	350	64	8100	670	59000	2054
3/28/2001	<5.7	36	140	12000	790	210000	1100000	100	110	860	4700	720	1400	870	29000	2600	100000	6142
6/27/2001	<2.8	49	220	21000	1100	300000	1300000	170	660	690	2200	7800	2300	1600	290000	4000	110000	11391.50

I-TEF/89 = International Toxicity Equivalent Factors/1989

I-TEQ/89 = International Toxicity Equivalents (based on I-TEF/89)

= Exceedance of I-TEQ/89

= Results are above detection limit but below quantitation limit

WEISENBERGER TIE & LUMBER COMPANY

Dioxin/Furan Analytical Results - DMW-2

Date Sampled	Compound (pg/l)																I-TEQ/89 2,3,7,8-TCDD	
	2,3,7,8-TCDD	1,2,3,7,8-PeCDD	1,2,3,4,7,8-HxCDD	1,2,3,6,7,8-HxCDD	1,2,3,7,8,9-HxCDD	1,2,3,4,6,7,8-HpCDD	1,2,3,4,6,7,8,9-OCDD	2,3,7,8-TCDF	1,2,3,7,8-PeCDF	2,3,4,7,8-PeCDF	1,2,3,4,7,8-HxCDF	1,2,3,6,7,8-HxCDF	2,3,4,6,7,8-HxCDF	1,2,3,7,8,9-HxCDF	1,2,3,4,6,7,8-HpCDF	1,2,3,4,7,8,9-HpCDF		
I-TEF/89	1.00	0.50	0.10	0.10	0.10	0.01	0.001	0.10	0.05	0.50	0.10	0.1	0.1	0.1	0.01	0.01	0.001	3.00
3/6/2000	<3.6	<3.6	<4.8	<4.4	<5.3	80	660	<3.0	<4.4	<3.2	<5.3	<4.1	<5.3	<2.6	10	<4.0	65	1.63
6/30/2000	<3.3	9.5	8.2	17	10	260	2400	<2.9	<5.5	<4.7	<11	<3.0	<2.6	<2.2	<12	<11	200	13.47
12/27/2000	NO SAMPLE																	
3/28/2001	<4.5	<2.9	<3.3	<3.0	<1.3	22	290	<3.7	<1.6	<1.5	<1.9	<1.2	<2.2	<2.7	2.6	<2.8	11	0.55
6/27/2001	<8.9	<9.5	<9.5	<9.5	<9.5	150	2300	<5.3	<9.5	<9.5	<9.5	<9.5	<9.5	<11	170	<27	280	5.78

I-TEF/89 = International Toxicity Equivalent Factors/1989

I-TEQ/89 = International Toxicity Equivalents (based on I-TEF/89)

= Exceedance of I-TEQ/89

= Results are above detection limit but below quantitation limit

WEISENBERGER TIE & LUMBER COMPANY

Dioxin/Furan Analytical Results - DMW-4

Date Sampled	Compound (pg/l)																I-TEQ/89 2,3,7,8-TCDD	
	2,3,7,8-TCDD	1,2,3,7,8-PeCDD	1,2,3,4,7,8-HxCDD	1,2,3,6,7,8-HxCDD	1,2,3,7,8,9-HxCDD	1,2,3,4,6,7,8-HpCDD	1,2,3,4,6,7,8,9-OCDD	2,3,7,8-TCDF	1,2,3,7,8-PeCDF	2,3,4,7,8-PeCDF	1,2,3,4,7,8-HxCDF	1,2,3,6,7,8-HxCDF	2,3,4,6,7,8-HxCDF	1,2,3,7,8,9-HxCDF	1,2,3,4,6,7,8-HpCDF	1,2,3,4,7,8,9-HpCDF		1,2,3,4,6,7,8,9-OCDF
I-TEF/89	1.00	0.50	0.10	0.10	0.10	0.01	0.001	0.10	0.05	0.50	0.10	0.1	0.1	0.1	0.01	0.01	0.001	3.00
8/20/1992	0	0	7.3	65.5	13.8	1950	13990	2.30	2.60	3.1	14.2	3.9	7.9	0	178	21	1340	49.99
6/30/2000	<4.8	21	18	140	40	1600	15000	<3.3	<19	<30	<8.7	<21	29	14	190	51	1300	69.31
9/27/2000	<11	<10	33	270	55	4800	33000	<13	<10	<8.2	38	16	49	25	500	55	2800	138
12/27/2000	NO SAMPLE																	
3/28/2001	<3.0	14	41	280	42	4000	31000	<4.4	<16	23	42	42	47	19	580	52	4000	151.12
06/27/2001	<7.9	<31	<36	64	<37	810	6700	<8.8	<22	<10	<21	<24	<24	<37	760	<17	550	29.35

I-TEF/89 = International Toxicity Equivalent Factors/1989

I-TEQ/89 = International Toxicity Equivalents (based on I-TEF/89)

= Exceedance of I-TEQ/89

= Results are above detection limit but below quantitation limit

WEISENBERGER TIE & LUMBER COMPANY

Dioxin/Furan Analytical Results - DMW-5

Date Sampled	Compound (pg/l)																I-TEQ/89 2,3,7,8-TCDD	
	2,3,7,8-TCDD	1,2,3,7,8-PeCDD	1,2,3,4,7,8-HxCDD	1,2,3,6,7,8-HxCDD	1,2,3,7,8,9-HxCDD	1,2,3,4,6,7,8-HpCDD	1,2,3,4,6,7,8,9-OCDD	2,3,7,8-TCDF	1,2,3,7,8-PeCDF	2,3,4,7,8-PeCDF	1,2,3,4,7,8-HxCDF	1,2,3,6,7,8-HxCDF	2,3,4,6,7,8-HxCDF	1,2,3,7,8,9-HxCDF	1,2,3,4,6,7,8-HpCDF	1,2,3,4,7,8,9-HpCDF		
I-TEF/89	1.00	0.50	0.10	0.10	0.10	0.01	0.001	0.10	0.05	0.50	0.10	0.1	0.1	0.1	0.01	0.01	0.001	3.00
1/19/1993	0	0	0	5.7	0	123	907	0	0	0	0	0	5.4	0	14.2	0	41.1	3.43
6/30/2000	<2.3	<1.6	<2.9	<2.4	<2.9	14	100	<1.1	<1.3	<1.3	<1.9	<2.1	<1.7	<1.8	<5.1	<10	11	0.25
9/27/2000	<18	<13	<20	<15	<14	290	2400	<11	<10	<7.9	<12	6.3	<18	<18	46	<26	200	6.59
3/28/2001	NO SAMPLE																	
6/27/2001	<4.7	<32	<36	<30	<41	290	1600	<4.7	<19	<12	<9.6	<9.4	<13	<29	190	<26	130	6.53

I-TEF/89 = International Toxicity Equivalent Factors/1989

I-TEQ/89 = International Toxicity Equivalents (based on I-TEF/89)

= Exceedance of I-TEQ/89

= Results are above detection limit but below quantitation limit

WEISENBERGER TIE & LUMBER COMPANY

Dioxin/Furan Analytical Results - DMW-6A

Date Sampled	Compound (pg/l)																I-TEQ/89 2,3,7,8-TCDD	
	2,3,7,8-TCDD	1,2,3,7,8-PeCDD	1,2,3,4,7,8-HxCDD	1,2,3,6,7,8-HxCDD	1,2,3,7,8,9-HxCDD	1,2,3,4,6,7,8-HpCDD	1,2,3,4,6,7,8,9-OCDD	2,3,7,8-TCDF	1,2,3,7,8-PeCDF	2,3,4,7,8-PeCDF	1,2,3,4,7,8-HxCDF	1,2,3,6,7,8-HxCDF	2,3,4,6,7,8-HxCDF	1,2,3,7,8,9-HxCDF	1,2,3,4,6,7,8-HpCDF	1,2,3,4,7,8,9-HpCDF		1,2,3,4,6,7,8,9-OCDF
I-TEF/89	1.00	0.50	0.10	0.10	0.10	0.01	0.001	0.10	0.05	0.50	0.10	0.1	0.1	0.1	0.01	0.01	0.001	3.00
8/20/1992	0	0	0	5.80	0	250.00	3170.000	0	0	0	0	0	3.9	0	17.90	2.90	304.000	7.15
6/30/2000	<5.2	15	13	38	16	970	13000	<2.9	<3.1	<2.2	<4.5	<13	<8.9	<5.1	100	<37	1200	39.10
12/27/2000	NO SAMPLE																	
3/28/2001	NO SAMPLE																	
6/27/2001	<7.4	<25	<33	<43	<36	90	580	<4.8	<19	<20	<21	<20	<15	<31	35	<35	38	1.87

I-TEF/89 = International Toxicity Equivalent Factors/1989

I-TEQ/89 = International Toxicity Equivalents (based on I-TEF/89)

= Exceedance of I-TEQ/89

= Results are above detection limit but below quantitation limit

WEISENBERGER TIE & LUMBER COMPANY

Dioxin/Furan Analytical Results - DMW-7

Date Sampled	Compound (pg/l)																I-TEQ/89 2,3,7,8-TCDD	
	2,3,7,8-TCDD	1,2,3,7,8-PeCDD	1,2,3,4,7,8-HxCDD	1,2,3,6,7,8-HxCDD	1,2,3,7,8,9-HxCDD	1,2,3,4,6,7,8-HpCDD	1,2,3,4,6,7,8,9-OCDD	2,3,7,8-TCDF	1,2,3,7,8-PeCDF	2,3,4,7,8-PeCDF	1,2,3,4,7,8-HxCDF	1,2,3,6,7,8-HxCDF	2,3,4,6,7,8-HxCDF	1,2,3,7,8,9-HxCDF	1,2,3,4,6,7,8-HpCDF	1,2,3,4,7,8,9-HpCDF		1,2,3,4,6,7,8,9-OCDF
I-TEF/89	1.00	0.50	0.10	0.10	0.10	0.01	0.001	0.10	0.05	0.50	0.10	0.1	0.1	0.1	0.01	0.01	0.001	3.00
8/20/1992	0	0	0	0	0	22.7	312	0	0	0	0	0	0	0	0	0	0	0.54
12/27/2000	<6.5	<5.5	<3.5	<2.9	<3.7	18	250	<5.4	<3.7	<3.7	18	<1.5	<2.3	<3.2	<1.4	<2.1	21	2.25
3/28/2001	<4.8	<1.8	<3.7	<3.6	<3.3	4.9	38	<2.9	<3.5	<2.4	<1.4	<1.5	<1.0	<1.9	<1.9	<1.6	<4.8	0.09
6/27/2001	<5.6	<45	<48	<42	<35	<46	<63	<4.9	<23	<17	<27	<22	<18	<16	<24	<34	<31	

I-TEF/89 = International Toxicity Equivalent Factors/1989

I-TEQ/89 = International Toxicity Equivalents (based on I-TEF/89)

= Exceedance of I-TEQ/89

= Results are above detection limit but below quantitation limit

WEISENBERGER TIE & LUMBER COMPANY

Dioxin/Furan Analytical Results - DMW-8

Date Sampled	Compound (pg/l)																	I-TEQ/89 2,3,7,8-TCDD
	2,3,7,8-TCDD	1,2,3,7,8-PeCDD	1,2,3,4,7,8-HxCDD	1,2,3,6,7,8-HxCDD	1,2,3,7,8,9-HxCDD	1,2,3,4,6,7,8-HpCDD	1,2,3,4,6,7,8,9-OCDD	2,3,7,8-TCDF	1,2,3,7,8-PeCDF	2,3,4,7,8-PeCDF	1,2,3,4,7,8-HxCDF	1,2,3,6,7,8-HxCDF	2,3,4,6,7,8-HxCDF	1,2,3,7,8,9-HxCDF	1,2,3,4,6,7,8-HpCDF	1,2,3,4,7,8,9-HpCDF	1,2,3,4,6,7,8,9-OCDF	
I-TEF/89	1.00	0.50	0.10	0.10	0.10	0.01	0.001	0.10	0.05	0.50	0.10	0.1	0.1	0.1	0.01	0.01	0.001	3.00
8/20/1992	0	0	0	0	0	152.00	2750.000	0	0	0	0	0	12.6	0	0	0	58.2	5.59
12/27/2000	<6.9	<3.1	<3.3	<3.1	<2.4	7.1	49	<4.7	<3.6	<1.7	<2.2	<2.3	<1.4	<1.7	5	<1.7	9.1	0.18
3/28/2001	<7.6	<5.5	<3.8	<2.4	<1.3	4.0	19	<5.7	<3.7	<2.2	<1.5	<2.5	<1.8	<0.74	<1.1	<1.7	4.0	0.06
6/27/2001	<8.4	<28	<32	<24	<33	<33	<42	<4.9	<20	<18	<22	<19	<16	<21	<30	<34	<33	

I-TEF/89 = International Toxicity Equivalent Factors/1989

I-TEQ/89 = International Toxicity Equivalents (based on I-TEF/89)

= Exceedance of I-TEQ/89

= Results are above detection limit but below quantitation limit

WEISENBERGER TIE & LUMBER COMPANY

Dioxin/Furan Analytical Results - DMW-10

Date Sampled	Compound (pg/l)																I-TEQ/89 2,3,7,8-TCDD	
	2,3,7,8-TCDD	1,2,3,7,8-PeCDD	1,2,3,4,7,8-HxCDD	1,2,3,6,7,8-HxCDD	1,2,3,7,8,9-HxCDD	1,2,3,4,6,7,8-HpCDD	1,2,3,4,6,7,8,9-OCDD	2,3,7,8-TCDF	1,2,3,7,8-PeCDF	2,3,4,7,8-PeCDF	1,2,3,4,7,8-HxCDF	1,2,3,6,7,8-HxCDF	2,3,4,6,7,8-HxCDF	1,2,3,7,8,9-HxCDF	1,2,3,4,6,7,8-HpCDF	1,2,3,4,7,8,9-HpCDF		1,2,3,4,6,7,8,9-OCDF
I-TEF/89	1.00	0.50	0.10	0.10	0.10	0.01	0.001	0.10	0.05	0.50	0.10	0.1	0.1	0.1	0.01	0.01	0.001	3.00
1/19/1993	0	0	0	0	0	28.7	449	0	0	0	0	0	3.1	0	0	0	6.4	1.05

I-TEF/89 = International Toxicity Equivalent Factors/1989

I-TEQ/89 = International Toxicity Equivalents (based on I-TEF/89)

= Exceedance of I-TEQ/89

= Results are above detection limit but below quantitation limit

WEISENBERGER TIE & LUMBER COMPANY

Dioxin/Furan Analytical Results - DMW-13

Date Sampled	Compound (pg/l)																I-TEQ/89 2,3,7,8-TCDD	
	2,3,7,8-TCDD	1,2,3,7,8-PeCDD	1,2,3,4,7,8-HxCDD	1,2,3,6,7,8-HxCDD	1,2,3,7,8,9-HxCDD	1,2,3,4,6,7,8-HpCDD	1,2,3,4,6,7,8,9-OCDD	2,3,7,8-TCDF	1,2,3,7,8-PeCDF	2,3,4,7,8-PeCDF	1,2,3,4,7,8-HxCDF	1,2,3,6,7,8-HxCDF	2,3,4,6,7,8-HxCDF	1,2,3,7,8,9-HxCDF	1,2,3,4,6,7,8-HpCDF	1,2,3,4,7,8,9-HpCDF		
I-TEF/89	1.00	0.50	0.10	0.10	0.10	0.01	0.001	0.10	0.05	0.50	0.10	0.1	0.1	0.1	0.01	0.01	0.001	3.00
12/23/1999	<6.6	<5.6	<6.7	<7.1	<6.3	<6.4	26	<3.2	<4.7	<3.6	<2.9	<2.8	<4.5	<4.4	<5.2	<4.1	<6.2	0.03
3/6/2000	<12	<12	<13	<12	<11	<12	44	<8.0	<8.8	<7.3	<7.0	<6.3	<6.4	<16	<6.3	<15	<21	0.04

I-TEF/89 = International Toxicity Equivalent Factors/1989

I-TEQ/89 = International Toxicity Equivalents (based on I-TEF/89)

= Exceedance of I-TEQ/89

= Results are above detection limit but below quantitation limit

WEISENBERGER TIE & LUMBER COMPANY

Dioxin/Furan Analytical Results - DPZ-1

Date Sampled	Compound (pg/l)																	I-TEQ/89 2,3,7,8-TCDD
	2,3,7,8-TCDD	1,2,3,7,8-PeCDD	1,2,3,4,7,8-HxCDD	1,2,3,6,7,8-HxCDD	1,2,3,7,8,9-HxCDD	1,2,3,4,6,7,8-HpCDD	1,2,3,4,6,7,8,9-OCDD	2,3,7,8-TCDF	1,2,3,7,8-PeCDF	2,3,4,7,8-PeCDF	1,2,3,4,7,8-HxCDF	1,2,3,6,7,8-HxCDF	2,3,4,6,7,8-HxCDF	1,2,3,7,8,9-HxCDF	1,2,3,4,6,7,8-HpCDF	1,2,3,4,7,8,9-HpCDF	1,2,3,4,6,7,8,9-OCDF	
I-TEF/89	1.00	0.50	0.10	0.10	0.10	0.01	0.001	0.10	0.05	0.50	0.10	0.1	0.1	0.1	0.01	0.01	0.001	3.00
8/20/1992	0	0	0	1750	159	54270	385020	21.2	83.6	73.1	245	96.5	150	0	4050	392	34250	1289.29
12/23/1999	<2.8	<3.3	<4.1	<4.0	<3.8	50	470	<2.3	<2.4	<1.6	<1.5	<1.1	<2.3	<1.9	6	<3.0	42	1.07
3/6/2000	<17	<17	<9.8	55	<11	1100	10000	14	<5.3	<14	<11	<12	<11	<16	86	12	970	29.85
6/30/2000	<3.0	<2.9	<3.4	56	5.0	1200	11000	<4.7	<3.0	<3.7	<4.0	<3.2	6.0	4.6	120	.13	890	32.38
9/27/2000	<5.2	<3.4	<6.9	74	<4.7	2100	22000	<3.3	<3.8	7.8	8.0	<3.2	11	<4.5	210	18	1700	60.18
12/27/2000	<2.7	<4.7	<1.6	260	12	5500	51000	<1.7	14	21	37	<2.0	33	17	<4.0	43	4200	157.73
03/28/2001	<4.0	<3.0	12	110	6.0	2600	25000	<3.2	<6.1	7.2	15	11	12	6.3	300	.20	2300	77.33
06/27/2001	<5.4	<13	<9.5	93	<22	1800	19000	<6.8	<13	<14	10	10	<11	<11	1300	29	1400	62.99

I-TEF/89 = International Toxicity Equivalent Factors/1989

I-TEQ/89 = International Toxicity Equivalents (based on I-TEF/89)

 = Exceedance of I-TEQ/89

 = Results are above detection limit but below quantitation limit

WEISENBERGER TIE & LUMBER COMPANY

Dioxin/Furan Analytical Results - DPZ-1a

Date Sampled	Compound (pg/l)																	I-TEQ/89 2,3,7,8-TCDD
	2,3,7,8-TCDD	1,2,3,7,8-PeCDD	1,2,3,4,7,8-HxCDD	1,2,3,6,7,8-HxCDD	1,2,3,7,8,9-HxCDD	1,2,3,4,6,7,8-HpCDD	1,2,3,4,6,7,8,9-OCDD	2,3,7,8-TCDF	1,2,3,7,8-PeCDF	2,3,4,7,8-PeCDF	1,2,3,4,7,8-HxCDF	1,2,3,6,7,8-HxCDF	2,3,4,6,7,8-HxCDF	1,2,3,7,8,9-HxCDF	1,2,3,4,6,7,8-HpCDF	1,2,3,4,7,8,9-HpCDF	1,2,3,4,6,7,8,9-OCDF	
I-TEF/89	1.00	0.50	0.10	0.10	0.10	0.01	0.001	0.10	0.05	0.50	0.10	0.1	0.1	0.1	0.01	0.01	0.001	3.00
8/20/1992	0	0	0	0	0	121	0	0	0	0	0	0	0	0	10.7	0	52.6	1.37

I-TEF/89 = International Toxicity Equivalent Factors/1989

I-TEQ/89 = International Toxicity Equivalents (based on I-TEF/89)

= Exceedance of I-TEQ/89

= Results are above detection limit but below quantitation limit

WEISENBERGER TIE & LUMBER COMPANY

Dioxin/Furan Analytical Results - DPZ-2

Date Sampled	Compound (pg/l)																I-TEQ/89 2,3,7,8-TCDD	
	2,3,7,8-TCDD	1,2,3,7,8-PeCDD	1,2,3,4,7,8-HxCDD	1,2,3,6,7,8-HxCDD	1,2,3,7,8,9-HxCDD	1,2,3,4,6,7,8-HpCDD	1,2,3,4,6,7,8,9-OCDD	2,3,7,8-TCDF	1,2,3,7,8-PeCDF	2,3,4,7,8-PeCDF	1,2,3,4,7,8-HxCDF	1,2,3,6,7,8-HxCDF	2,3,4,6,7,8-HxCDF	1,2,3,7,8,9-HxCDF	1,2,3,4,6,7,8-HpCDF	1,2,3,4,7,8,9-HpCDF		1,2,3,4,6,7,8,9-OCDF
I-TEF/89	1.00	0.50	0.10	0.10	0.10	0.01	0.001	0.10	0.05	0.50	0.10	0.1	0.1	0.1	0.01	0.01	0.001	3.00
12/23/1999	<4.3	<4.5	<5.9	<5.4	<5.4	<8.3	140	<2.8	<3.1	<4.3	<1.9	<2.1	<5.8	<8.9	<5.1	<8.1	25	0.17
3/6/2000	<3.0	<4.5	<5.1	<4.7	<4.8	<4.4	25	<2.6	<4.1	<3.1	<2.6	<1.9	<2.5	<3.9	<3.2	<2.6	<4.6	0.03

I-TEF/89 = International Toxicity Equivalent Factors/1989

I-TEQ/89 = International Toxicity Equivalents (based on I-TEF/89)

= Exceedance of I-TEQ/89

= Results are above detection limit but below quantitation limit

WEISENBERGER TIE & LUMBER COMPANY

Dioxin/Furan Analytical Results - DPZ-3

Date Sampled	Compound (pg/l)																I-TEQ/89 2,3,7,8-TCDD	
	2,3,7,8-TCDD	1,2,3,7,8-PeCDD	1,2,3,4,7,8-HxCDD	1,2,3,6,7,8-HxCDD	1,2,3,7,8,9-HxCDD	1,2,3,4,6,7,8-HpCDD	1,2,3,4,6,7,8,9-OCDD	2,3,7,8-TCDF	1,2,3,7,8-PeCDF	2,3,4,7,8-PeCDF	1,2,3,4,7,8-HxCDF	1,2,3,6,7,8-HxCDF	2,3,4,6,7,8-HxCDF	1,2,3,7,8,9-HxCDF	1,2,3,4,6,7,8-HpCDF	1,2,3,4,7,8,9-HpCDF		1,2,3,4,6,7,8,9-OCDF
I-TEF/89	1.00	0.50	0.10	0.10	0.10	0.01	0.001	0.10	0.05	0.50	0.10	0.1	0.1	0.1	0.01	0.01	0.001	3.00
8/20/1992	0	0	0	0	0	45.9	633	0	0	0	2.1	0	4.2	0	3.8	0	18.1	1.78
12/27/2000	<7.1	<3.6	<3.1	<2.6	<3.4	21	450	<5.9	<3.3	<2.7	<3.4	<3.8	<2.0	<2.5	<1.8	<2.9	29	0.69
3/28/2001	<5.0	<3.3	<1.2	<2.3	<1.9	20	530	<3.3	<2.0	<1.6	<2.2	<2.3	<1.9	<1.2	7.8	<2.1	19	0.83
6/27/2001	<6.7	<17	<15	<17	<15	<27	200	<5.0	<14	<9.6	<10	<11	<9.7	<13	<13	19	<19	0.39

I-TEF/89 = International Toxicity Equivalent Factors/1989

I-TEQ/89 = International Toxicity Equivalents (based on I-TEF/89)

= Exceedance of I-TEQ/89

= Results are above detection limit but below quantitation limit

WEISENBERGER TIE & LUMBER COMPANY

Dioxin/Furan Analytical Results - DPZ-4

Date Sampled	Compound (pg/l)																I-TEQ/89 2,3,7,8-TCDD	
	2,3,7,8-TCDD	1,2,3,7,8-PeCDD	1,2,3,4,7,8-HxCDD	1,2,3,6,7,8-HxCDD	1,2,3,7,8,9-HxCDD	1,2,3,4,6,7,8-HpCDD	1,2,3,4,6,7,8,9-OCDD	2,3,7,8-TCDF	1,2,3,7,8-PeCDF	2,3,4,7,8-PeCDF	1,2,3,4,7,8-HxCDF	1,2,3,6,7,8-HxCDF	2,3,4,6,7,8-HxCDF	1,2,3,7,8,9-HxCDF	1,2,3,4,6,7,8-HpCDF	1,2,3,4,7,8,9-HpCDF		1,2,3,4,6,7,8,9-OCDF
I-TEF/89	1.00	0.50	0.10	0.10	0.10	0.01	0.001	0.10	0.05	0.50	0.10	0.1	0.1	0.1	0.01	0.01	0.001	3.00
12/23/1999	<18	<14	<12	<16	<14	23	190	<6.9	<10	<8.6	<12	<12	<6.1	<11	<26	<22	<27	0.42
3/6/2000	<9.2	<6.4	<7.3	<7.1	<8.6	11	56	<4.9	<8.7	<5.3	<2.7	<5.4	<6.2	<5.8	<5.9	<8.3	11	0.18

I-TEF/89 = International Toxicity Equivalent Factors/1989

I-TEQ/89 = International Toxicity Equivalents (based on I-TEF/89)

= Exceedance of I-TEQ/89

= Results are above detection limit but below quantitation limit

WEISENBERGER TIE & LUMBER COMPANY

Dioxin/Furan Analytical Results - DPZ-5

Date Sampled	Compound (pg/l)															I-TEQ/89 2,3,7,8-TCDD		
	2,3,7,8-TCDD	1,2,3,7,8-PeCDD	1,2,3,4,7,8-HxCDD	1,2,3,6,7,8-HxCDD	1,2,3,7,8,9-HxCDD	1,2,3,4,6,7,8-HpCDD	1,2,3,4,6,7,8,9-OCDD	2,3,7,8-TCDF	1,2,3,7,8-PeCDF	2,3,4,7,8-PeCDF	1,2,3,4,7,8-HxCDF	1,2,3,6,7,8-HxCDF	2,3,4,6,7,8-HxCDF	1,2,3,7,8,9-HxCDF	1,2,3,4,6,7,8-HpCDF		1,2,3,4,7,8,9-HpCDF	1,2,3,4,6,7,8,9-OCDF
I-TEF/89	1.00	0.50	0.10	0.10	0.10	0.01	0.001	0.10	0.05	0.50	0.10	0.1	0.1	0.1	0.01	0.01	0.001	3.00
8/20/1992	0	0	0	0	0	23.2	309	0	0	0	0	0	0	0	0	0	0	0.54

I-TEF/89 = International Toxicity Equivalent Factors/1989

I-TEQ/89 = International Toxicity Equivalents (based on I-TEF/89)

= Exceedance of I-TEQ/89

= Results are above detection limit but below quantitation limit

WEISENBERGER TIE & LUMBER COMPANY

Dioxin/Furan Analytical Results - DPZ-6

Date Sampled	Compound (pg/l)																I-TEQ/89 2,3,7,8-TCDD	
	2,3,7,8-TCDD	1,2,3,7,8-PeCDD	1,2,3,4,7,8-HxCDD	1,2,3,6,7,8-HxCDD	1,2,3,7,8,9-HxCDD	1,2,3,4,6,7,8-HpCDD	1,2,3,4,6,7,8,9-OCDD	2,3,7,8-TCDF	1,2,3,7,8-PeCDF	2,3,4,7,8-PeCDF	1,2,3,4,7,8-HxCDF	1,2,3,6,7,8-HxCDF	2,3,4,6,7,8-HxCDF	1,2,3,7,8,9-HxCDF	1,2,3,4,6,7,8-HpCDF	1,2,3,4,7,8,9-HpCDF		1,2,3,4,6,7,8,9-OCDF
I-TEF/89	1.00	0.50	0.10	0.10	0.10	0.01	0.001	0.10	0.05	0.50	0.10	0.1	0.1	0.1	0.01	0.01	0.001	3.00
12/23/1999	<2.2	<3.1	<3.0	<3.9	<3.7	<3.3	110	<1.9	<1.7	<1.8	<2.7	<1.8	<1.9	<2.1	3.4	<3.1	12	0.16
12/27/2000	NO SAMPLE																	
3/28/2001	<5.1	<1.4	<4.5	<3.8	<2.2	20	170	<3.8	<2.7	<1.9	<1.6	<1.8	<3.1	<2.0	5.7	<1.4	16	0.44
6/27/2001	<1.9	<13	<25	13	<27	340	3900	<1.9	<10	<9.4	<15	<15	<15	<13	380	<24	320	12.72

I-TEF/89 = International Toxicity Equivalent Factors/1989

I-TEQ/89 = International Toxicity Equivalents (based on I-TEF/89)

= Exceedance of I-TEQ/89


= Results are above detection limit but below quantitation limit

MONITORING WELLS
PVOC ANALYSIS

WEISENBERGER TIE & LUMBER COMPANY
PVOC Analytical Results

MONITORING WELL MW-1								
Date Sampled	Benzene (ug/L)	Toluene (ug/L)	Ethylbenzene (ug/L)	Xylenes (ug/L)	1,2,4 TMB (ug/L)	1,3,5 TMB (ug/L)	MTBE (ug/L)	comments
WDNR ES	5	343	700	620			60	
06/03/92	<1.0	<1.0	<1.0	<3.0	<1.0	<1.0	NA	
08/20/92	<1.0	<1.0	<1.0	<3.0	<1.0	<1.0	NA	
12/20/94	-	-	-	-	-	-	-	NS
03/11/98	-	-	-	-	-	-	-	NS
06/23/98	-	-	-	-	-	-	-	NS

MONITORING WELL MW-2								
Date Sampled	Benzene (ug/L)	Toluene (ug/L)	Ethylbenzene (ug/L)	Xylenes (ug/L)	1,2,4 TMB (ug/L)	1,3,5 TMB (ug/L)	MTBE (ug/L)	comments
WDNR ES	5	343	700	620			60	
06/03/92	<1.0	<1.0	<1.0	<3.0	<1.0	<1.0	NA	
08/20/92	<1.0	<1.0	<1.0	<3.0	<1.0	<1.0	NA	
12/20/94	<1.0	1	<1.0	<3.0	<1.0	<1.0	NA	
03/11/98	-	-	-	-	-	-	-	NS
06/23/98	-	-	-	-	-	-	-	NS

 = ES exceedance

WEISENBERGER TIE & LUMBER COMPANY
PVOC Analytical Results

MONITORING WELL								MW-3
Date Sampled	Benzene (ug/L)	Toluene (ug/L)	Ethylbenzene (ug/L)	Xylenes (ug/L)	1,2,4 TMB (ug/L)	1,3,5 TMB (ug/L)	MTBE (ug/L)	comments
WDNR ES	5	343	700	620			60	
06/03/92	7.9	21	21	150	85	20	NA	
08/20/92	<10	16	15	150	84	15	NA	
12/20/94	<10	<10	15	120	90	<10	NA	
03/11/98	3.7	1.9	14	85	80	14	<1.3	
06/24/98	2.9	1.6	11	71	71	15	<0.8	

MW-3

Date Sampled	Benzene µg/L	Toluene µg/L	Ethylbenzene µg/L	Xylenes µg/L	Trimethybenzenes µg/L	MTBE µg/L
WDNR ES	5	1000	700	10000	480	60
9/15/98	4.1	2.1	17	104	123	<0.92
12/2/98	4.2	2.1	17	109	131	<0.92
3/30/99	3.7	1.7	13	84	100	<0.92
6/10/99	5.0	2.8	17	105	123	<0.92
9/20/99	5.4	2.9	17	106	136	<0.92
12/3/99	4.1	2.1	13	87	105	<0.92
6/30/00	4.2	2.2	13	77	96	<0.92
9/27/00	4.1	2.3	14	95	85.1	<0.92
12/27/00	<5.0	<6.0	12	77	126	<9.2
3/28/01	2.9	1.4	9.1	54	69	<0.091
6/27/01	3.6	2.0	11	69	90	<0.091

= ES exceedance

WEISENBERGER TIE & LUMBER COMPANY
PVOC Analytical Results

MONITORING WELL		MW-5						
Date Sampled	Benzene (ug/L)	Toluene (ug/L)	Ethylbenzene (ug/L)	Xylenes (ug/L)	1,2,4 TMB (ug/L)	1,3,5 TMB (ug/L)	MTBE (ug/L)	comments
WDNR ES	5	343	700	620			60	
06/03/92	10	5.5	26	400	340	75	NA	
08/20/92	<5.0	<5.0	5.7	100	100	20	NA	
12/20/94	<5.0	<5.0	5.4	47	94	17	NA	
03/11/98	<0.13	0.20	<0.22	8.4	11	1.7	0.7	
06/24/98	0.23	<0.20	<0.22	20	25	2.6	<0.16	

MW-5

Date Sampled	Benzene µg/L	Toluene µg/L	Ethylbenzene µg/L	Xylenes µg/L	Trimethybenzenes µg/L	MTBE µg/L
WDNR ES	5	1000	700	10000	480	60
9/15/98	<0.50	<0.60	<0.60	34.7	49.7	<0.92
12/2/98	<0.50	<0.60	<0.60	38	52.6	<0.92
3/30/99	<0.50	<0.60	<0.60	33.6	40.5	<0.92
6/10/99	<0.50	<0.60	<0.60	38.7	50.3	<0.92
9/20/99	<0.50	<0.60	<0.60	36.9	56.4	<0.92
12/3/99	<0.50	<0.60	<0.60	34	43.6	<0.92

= ES exceedance

WEISENBERGER TIE & LUMBER COMPANY
PVOC Analytical Results

MONITORING WELL		MW-6						
Date Sampled	Benzene (ug/L)	Toluene (ug/L)	Ethylbenzene (ug/L)	Xylenes (ug/L)	1,2,4 TMB (ug/L)	1,3,5 TMB (ug/L)	MTBE (ug/L)	comments
WDNR ES	5	343	700	620			60	
06/03/92	<1.0	<1.0	<1.0	<3.0	<1.0	<1.0	NA	
08/20/92	2.1	10	2.4	15	5	1.2	NA	
12/20/94	<1.0	<1.0	<1.0	<3.0	<1.0	<1.0	NA	
03/11/98	-	-	-	-	-	-	-	NS
06/23/98	-	-	-	-	-	-	-	NS

MW-6

Date Sampled	Benzene µg/L	Toluene µg/L	Ethylbenzene µg/L	Xylenes µg/L	Trimethybenzenes µg/L	MTBE µg/L
WDNR ES	5	1000	700	10000	480	60
3/6/00	<0.5	<0.6	<0.6	<1.7	<1.7	<0.92

= ES exceedance

WEISENBERGER TIE & LUMBER COMPANY
PVOC Analytical Results

MONITORING WELL								MW-7
Date Sampled	Benzene (ug/L)	Toluene (ug/L)	Ethylbenzene (ug/L)	Xylenes (ug/L)	1,2,4 TMB (ug/L)	1,3,5 TMB (ug/L)	MTBE (ug/L)	comments
WDNR ES	5	143	700	620			60	
06/03/92	<5.0	<5.0	<5.0	16	55	16	NA	
08/20/92	<5.0	<5.0	<5.0	14	50	12	NA	
12/20/94	<5.0	<5.0	<5.0	15	53	12	NA	
03/11/98	-	-	-	-	-	-	-	NS
06/23/98	-	-	-	-	-	-	-	NS

MW-7

Date Sampled	Benzene µg/L	Toluene µg/L	Ethylbenzene µg/L	Xylenes µg/L	Trimethybenzenes µg/L	MTBE µg/L
WDNR ES	5	1000	700	10000	480	60
3/6/00	<0.5	<0.6	1.1	7.6	18.7	<0.92

= ES exceedance

WEISENBERGER TIE & LUMBER COMPANY
PVOC Analytical Results

MONITORING WELL		MW-10						
Date Sampled	Benzene (ug/L)	Toluene (ug/L)	Ethylbenzene (ug/L)	Xylenes (ug/L)	1,2,4 TMB (ug/L)	1,3,5 TMB (ug/L)	MTBE (ug/L)	comments
WDNR ES	5	343	700	620			60	
06/03/92	<5.0	<5.0	<5.0	<15	46	10	NA	
08/20/92	<1.0	<1.0	<1.0	4.6	28	3.9	NA	
12/20/94	<1.0	<1.0	<1.0	<3.0	17	5.9	NA	
03/11/98	-	-	-	-	-	-	-	NS
06/23/98	-	-	-	-	-	-	-	NS

MW-10

Date Sampled	Benzene µg/L	Toluene µg/L	Ethylbenzene µg/L	Xylenes µg/L	Trimethylbenzenes µg/L	MTBE µg/L
WDNR ES	5	1000	700	10000	480	60
3/6/00	<0.5	<0.6	<0.6	<1.7	10.8	<0.92


= ES exceedance

WEISENBERGER TIE & LUMBER COMPANY
PVOC Analytical Results

MONITORING WELL DMW-1								
Date Sampled	Benzene (ug/L)	Toluene (ug/L)	Ethylbenzene (ug/L)	Xylenes (ug/L)	1,2,4 TMB (ug/L)	1,3,5 TMB (ug/L)	MTBE (ug/L)	comments
WDNR ES	5	343	700	620			60	
06/03/92	<5.0	<5.0	<5.0	<15	12	17	NA	
08/20/92	<5.0	21	13	113	72	17	NA	
12/20/94	-	-	-	-	-	-	-	NS
03/11/98	<0.13	2.1	2.5	21	17	18	0.27	
06/24/98	<0.13	1.1	1.6	14	18	14	<3.5	

DMW-1

Date Sampled	Benzene µg/L	Toluene µg/L	Ethylbenzene µg/L	Xylenes µg/L	Trimethybenzenes µg/L	MTBE µg/L
WDNR ES	5	1000	700	10000	480	60
9/15/98	<0.50	3.3	4.3	47	85	<0.92
12/2/98	<0.50	3.5	4.6	49	87	<0.92
3/30/99	<0.50	3.8	4.6	47	82	<0.92
6/10/99	<0.50	0.97	1.1	10.8	34	<0.92
9/20/99	<0.50	1.0	1.3	11.9	37	<0.92
12/3/99	<0.50	3.0	3.7	38	73	<0.92

 = ES exceedance

WEISENBERGER TIE & LUMBER COMPANY
PVOC Analytical Results

MONITORING WELL								DMW-2
Date Sampled	Benzene (ug/L)	Toluene (ug/L)	Ethylbenzene (ug/L)	Xylenes (ug/L)	1,2,4 TMB (ug/L)	1,3,5 TMB (ug/L)	MTBE (ug/L)	comments
WDNR ES	5	343	700	620			60	
06/03/92	<1.0	<1.0	<1.0	<3.0	<1.0	<1.0	NA	
08/20/92	-	-	-	-	-	-	-	NS
12/20/94	-	-	-	-	-	-	-	NS
03/11/98	<0.13	<0.20	<0.22	<0.23	<0.22	<0.29	<0.16	
06/24/98	<0.13	<0.20	<0.22	<0.23	<0.22	<0.29	<0.16	

DMW-2

Date Sampled	Benzene µg/L	Toluene µg/L	Ethylbenzene µg/L	Xylenes µg/L	Trimethybenzenes µg/L	MTBE µg/L
WDNR ES	5	1000	700	10000	480	60
9/15/98	Not Sampled					
12/2/98	Not Sampled					
3/30/99	<0.50	<0.60	<0.60	<1.7	<1.7	<0.92
6/10/99	<0.50	<0.60	<0.60	<1.7	<1.7	<0.92
9/20/99	Not Sampled					
12/3/99	Not Sampled					

= ES exceedance

WEISENBERGER TIE & LUMBER COMPANY
PVOC Analytical Results

MONITORING WELL		DMW-3						
Date Sampled	Benzene (ug/L)	Toluene (ug/L)	Ethylbenzene (ug/L)	Xylenes (ug/L)	1,2,4 TMB (ug/L)	1,3,5 TMB (ug/L)	MTBE (ug/L)	comments
WDNR ES	5	343	700	620			60	
06/03/92	<1.0	<1.0	<1.0	<3.0	<1.0	<1.0	NA	
08/20/92	-	-	-	-	-	-	-	NS
12/20/94	-	-	-	-	-	-	-	NS
03/11/98	-	-	-	-	-	-	-	NS
06/24/98	-	-	-	-	-	-	-	NS

DMW-3

Date Sampled	Benzene µg/L	Toluene µg/L	Ethylbenzene µg/L	Xylenes µg/L	Trimethybenzenes µg/L	MTBE µg/L
WDNR ES	5	1000	700	10000	480	60
3/6/00	Not Sampled					
6/30/00	Not Sampled					
9/27/00	Not Sampled					
12/27/00	Not Sampled					
3/28/01	Not Sampled					
6/27/01	Not Sampled					

= ES exceedance

WEISENBERGER TIE & LUMBER COMPANY
PVOC Analytical Results

MONITORING WELL DMW-4								
Date Sampled	Benzene (ug/L)	Toluene (ug/L)	Ethylbenzene (ug/L)	Xylenes (ug/L)	1,2,4 TMB (ug/L)	1,3,5 TMB (ug/L)	MTBE (ug/L)	comments
WDNR ES	5	343	700	620			60	
06/03/92	<1.0	1.5	1.4	18	19	10	NA	
08/20/92	<1.0	1.5	<1.0	16	17	7.9	NA	
12/20/94	-	-	-	-	-	-	-	NS
03/11/98	-	-	-	-	-	-	-	NS
06/24/98	-	-	-	-	-	-	-	NS

DMW-4

Date Sampled	Benzene µg/L	Toluene µg/L	Ethylbenzene µg/L	Xylenes µg/L	Trimethylbenzenes µg/L	MTBE µg/L
WDNR ES	5	1000	700	10000	480	60
3/6/00	Not Sampled					
6/30/00	<0.5	<0.6	<0.6	<1.7	<1.7	<0.92
9/27/00	<0.5	<0.6	<0.6	<1.7	<1.7	<0.92
12/27/00	Not Sampled					
3/28/00	<0.21	<0.22	<0.23	<0.44	<0.23	<0.091
6/27/01	<0.21	<0.22	<0.23	2.9	4.4	<0.091

= ES exceedance

WEISENBERGER TIE & LUMBER COMPANY
PVOC Analytical Results

MONITORING WELL		DMW-5						
Date Sampled	Benzene (ug/L)	Toluene (ug/L)	Ethylbenzene (ug/L)	Xylenes (ug/L)	1,2,4 TMB (ug/L)	1,3,5 TMB (ug/L)	MTBE (ug/L)	comments
WDNR ES	5	343	700	620			60	
06/03/92	<1.0	<1.0	<1.0	<3.0	<1.0	<1.0	NA	
08/20/92	<1.0	<1.0	<1.0	<3.0	<1.0	<1.0	NA	
12/20/94	<1.0	<1.0	<1.0	<3.0	<1.0	<1.0	NA	
03/11/98	-	-	-	-	-	-	-	NS
06/24/98	-	-	-	-	-	-	-	NS

DMW-5

Date Sampled	Benzene µg/L	Toluene µg/L	Ethylbenzene µg/L	Xylenes µg/L	Trimethybenzenes µg/L	MTBE µg/L
WDNR ES	5	1000	700	10000	480	60
3/6/00	Not Sampled					
6/30/00	<0.5	<0.6	<0.6	<1.7	<1.7	<0.92
9/27/00	<0.5	4.2	<0.6	<1.7	<1.7	<0.92
12/27/00	Not Sampled					
3/28/01	Not Sampled					
6/27/01	<0.21	<0.22	<0.23	<0.44	<0.23	<0.091

= ES exceedance

WEISENBERGER TIE & LUMBER COMPANY
PVOC Analytical Results

MONITORING WELL								DMW-6a
Date Sampled	Benzene (ug/L)	Toluene (ug/L)	Ethylbenzene (ug/L)	Xylenes (ug/L)	1,2,4 TMB (ug/L)	1,3,5 TMB (ug/L)	MTBE (ug/L)	comments
WDNR ES	5	343	700	620			60	
06/03/92	<1.0	<1.0	<1.0	<3.0	<1.0	<1.0	NA	
08/20/92	1.9	9.4	2	14	4.7	<1.0	NA	
12/20/94	<1.0	<1.0	<1.0	<3.0	<1.0	<1.0	NA	
03/11/98	-	-	-	-	-	-	-	NS
06/24/98	-	-	-	-	-	-	-	NS

DMW-6A

Date Sampled	Benzene µg/L	Toluene µg/L	Ethylbenzene µg/L	Xylenes µg/L	Trimethybenzenes µg/L	MTBE µg/L
WDNR ES	5	1000	700	10000	480	60
3/6/00	Not Sampled					
6/30/00	<0.5	<0.6	<0.6	<1.7	<1.7	<0.92
9/27/00	<0.5	<0.6	<0.6	<1.7	<1.7	<0.92
12/27/00	Not Sampled					
3/28/01	Not Sampled					
6/27/01	<0.21	<0.22	<0.23	<0.44	<0.23	<0.091

= ES exceedance

WEISENBERGER TIE & LUMBER COMPANY
PVOC Analytical Results

MONITORING WELL DMW-7								
Date Sampled	Benzene (ug/L)	Toluene (ug/L)	Ethylbenzene (ug/L)	Xylenes (ug/L)	1,2,4 TMB (ug/L)	1,3,5 TMB (ug/L)	MTBE (ug/L)	comments
WDNR ES	5	343	700	620			60	
06/03/92	-	-	-	-	-	-	-	NS
08/20/92	<1.0	<1.0	<1.0	<3.0	<1.0	<1.0	NA	
12/20/94	<1.0	<1.0	<1.0	<3.0	<1.0	<1.0	NA	
03/11/98	<0.13	<0.20	<0.22	<0.23	<0.22	<0.29	<0.16	
06/24/98	<0.13	<0.20	<0.22	<0.23	<0.22	<0.29	<0.16	

DMW-7

Date Sampled	Benzene µg/L	Toluene µg/L	Ethylbenzene µg/L	Xylenes µg/L	Trimethylbenzenes µg/L	MTBE µg/L
WDNR ES	5	1000	700	10000	480	60
9/15/98	<0.50	<0.60	<0.60	<1.7	<1.7	<0.92
12/2/98	<0.50	<0.60	<0.60	<1.7	<1.7	<0.92
3/30/99	<0.50	<0.60	<0.60	<1.7	<1.7	<0.92
6/10/99	<0.50	<0.60	<0.60	<1.7	<1.7	<0.92
9/20/99	<0.50	<0.60	<0.60	<1.7	<1.7	<0.92
12/3/99	<0.50	<0.60	<0.60	<1.7	<1.7	<0.92

= ES exceedance

WEISENBERGER TIE & LUMBER COMPANY
PVOC Analytical Results

MONITORING WELL								DMW-8
Date Sampled	Benzene (ug/L)	Toluene (ug/L)	Ethylbenzene (ug/L)	Xylenes (ug/L)	1,2,4 TMB (ug/L)	1,3,5 TMB (ug/L)	MTBE (ug/L)	comments
WDNR ES	5	343	700	620			60	
06/03/92	-	-	-	-	-	-	-	NS
08/20/92	<1.0	<1.0	<1.0	3.3	<1.0	<1.0	NA	
12/20/94	<1.0	<1.0	<1.0	<3.0	<1.0	<1.0	NA	
03/11/98	-	-	-	-	-	-	-	NS
06/24/98	-	-	-	-	-	-	-	NS

DMW-8

Date Sampled	Benzene µg/L	Toluene µg/L	Ethylbenzene µg/L	Xylenes µg/L	Trimethybenzenes µg/L	MTBE µg/L
WDNR ES	5	1000	700	10000	480	60
3/6/00	<0.50	<0.60	<0.60	<1.7	<1.7	<0.92

= ES exceedance

WEISENBERGER TIE & LUMBER COMPANY
PVOC Analytical Results

MONITORING WELL		DPZ-1						
Date Sampled	Benzene (ug/L)	Toluene (ug/L)	Ethylbenzene (ug/L)	Xylenes (ug/L)	1,2,4 TMB (ug/L)	1,3,5 TMB (ug/L)	MTBE (ug/L)	comments
WDNR ES	5	343	700	620			60	
06/03/92	93	670	68	300	900	300	NA	NS
08/20/92	1.6	20	<1.0	7.3	4	4.1	NA	NS
12/20/94	-	-	-	-	-	-	-	NS
03/11/98	0.25	3.30	<0.22	0.74	0.94	0.43	2	
06/24/98	0.31	2.40	<0.22	1.4	<0.22	<0.29	<0.16	

DPZ-1

Date Sampled	Benzene µg/L	Toluene µg/L	Ethylbenzene µg/L	Xylenes µg/L	Trimethylbenzenes µg/L	MTBE µg/L
WDNR ES	5	1000	700	10000	480	60
9/15/98	2.1	7.2	7.8	109	75	<0.92
12/2/98	2.2	10	12	131	115	<0.92
3/30/99	2.4	13	13	133	121	<0.92
6/10/99	2.6	6.0	14	143	130	<0.92
9/20/99	2.6	9.4	10	120	103	<0.92
12/3/99	2.5	2.7	14	139	120	<0.92
6/30/00	<5.0	<6.0	12	117	109	<9.2
9/27/00	1.8	15	7.5	119	67.8	<0.92
12/27/00	2.1	14	10	96	87	<0.091
3/28/01	2.0	3.4	10	70	91	<0.46
6/27/01	9.7	25	29	141	65	<1.8

= ES exceedance

WEISENBERGER TIE & LUMBER COMPANY
PVOC Analytical Results

MONITORING WELL								DMW-10
Date Sampled	Benzene (ug/L)	Toluene (ug/L)	Ethylbenzene (ug/L)	Xylenes (ug/L)	1,2,4 TMB (ug/L)	1,3,5 TMB (ug/L)	MTBE (ug/L)	comments
WDNR ES	5	343	700	620			60	
06/03/92	-	-	-	-	-	-	-	NS
08/20/92	-	-	-	-	-	-	-	NS
12/20/94	<1.0	<1.0	<1.0	<3.0	<1.0	<1.0	NA	
03/11/98	-	-	-	-	-	-	-	NS
06/24/98	-	-	-	-	-	-	-	NS

DMW-10

Date Sampled	Benzene µg/L	Toluene µg/L	Ethylbenzene µg/L	Xylenes µg/L	Trimethybenzenes µg/L	MTBE µg/L
WDNR ES	5	1000	700	10000	480	60
3/6/00	<0.50	<0.60	<0.60	<1.7	<1.7	<0.92
6/30/00	<0.50	<0.6	<0.6	<1.7	<1.7	<0.92
9/27/00	<0.5	<0.6	<0.6	<1.7	<1.7	<0.92
12/27/00	<0.5	<0.6	<0.6	<1.7	<1.7	<0.92
3/28/01	<0.21	<0.22	<0.23	<0.44	0.28	<0.091
6/27/01	<0.21	<0.22	<0.23	<0.44	<0.23	<0.091

= ES exceedance

WEISENBERGER TIE & LUMBER COMPANY
PVOC Analytical Results

MONITORING WELL		DPZ-1						
Date Sampled	Benzene (ug/L)	Toluene (ug/L)	Ethylbenzene (ug/L)	Xylenes (ug/L)	1,2,4 TMB (ug/L)	1,3,5 TMB (ug/L)	MTBE (ug/L)	comments
WDNR ES	5	343	700	620			60	
06/03/92	93	670	68	800	900	300	NA	NS
08/20/92	1.6	20	<1.0	7.3	4	4.1	NA	NS
12/20/94	-	-	-	-	-	-	-	NS
03/11/98	0.25	3.30	<0.22	0.74	0.94	0.43	2	
06/24/98	0.31	2.40	<0.22	1.4	<0.22	<0.29	<0.16	

DPZ-1

Date Sampled	Benzene µg/L	Toluene µg/L	Ethylbenzene µg/L	Xylenes µg/L	Trimethybenzenes µg/L	MTBE µg/L
WDNR ES	5	1000	700	10000	480	60
9/15/98	2.1	7.2	7.8	109	75	<0.92
12/2/98	2.2	10	12	131	115	<0.92
3/30/99	2.4	13	13	133	121	<0.92
6/10/99	2.6	6.0	14	143	130	<0.92
9/20/99	2.6	9.4	10	120	103	<0.92
12/3/99	2.5	2.7	14	139	120	<0.92
6/30/00	<5.0	<6.0	12	117	109	<9.2
9/27/00	1.8	15	7.5	119	67.8	<0.92
12/27/00	2.1	14	10	96	87	<0.091
3/28/01	2.0	3.4	10	70	91	<0.46
6/27/01	9.7	25	29	141	65	<1.8

= ES exceedance

WEISENBERGER TIE & LUMBER COMPANY
PVOC Analytical Results

MONITORING WELL DPZ-1a								
Date Sampled	Benzene (ug/L)	Toluene (ug/L)	Ethylbenzene (ug/L)	Xylenes (ug/L)	1,2,4 TMB (ug/L)	1,3,5 TMB (ug/L)	MTBE (ug/L)	comments
WDNR ES	5	343	700	620			60	
06/03/92	-	-	-	-	-	-	-	NS
08/20/92	<1.0	<1.0	<1.0	<3.0	<1.0	<1.0	NA	
12/20/94	1.7	2.1	<1.0	<3.0	<1.0	<1.0	NA	
03/11/98	-	-	-	-	-	-	-	NS
06/24/98	-	-	-	-	-	-	-	NS

DPZ-1a

Date Sampled	Benzene µg/L	Toluene µg/L	Ethylbenzene µg/L	Xylenes µg/L	Trimethybenzenes µg/L	MTBE µg/L
WDNR ES	5	1000	700	10000	480	60
3/6/00	<0.50	<0.60	<0.60	<1.7	<1.7	<0.92

= ES exceedance

WEISENBERGER TIE & LUMBER COMPANY
PVOC Analytical Results

MONITORING WELL DPZ-2								
Date Sampled	Benzene (ug/L)	Toluene (ug/L)	Ethylbenzene (ug/L)	Xylenes (ug/L)	1,2,4 TMB (ug/L)	1,3,5 TMB (ug/L)	MTBE (ug/L)	comments
WDNR ES	5	343	700	620			60	
06/03/92	<1.0	<1.0	<1.0	<3.0	<1.0	<1.0	NA	
08/20/92	<1.0	<1.0	<1.0	<3.0	<1.0	<1.0	NA	
12/20/94	<1.0	<1.0	<1.0	<3.0	<1.0	<1.0	NA	
03/11/98	<0.13	<0.20	<0.22	<0.23	<0.22	<0.29	<0.16	
06/24/98	<0.13	<0.20	<0.22	<0.23	<0.22	<0.29	<0.16	

DPZ-2

Date Sampled	Benzene µg/L	Toluene µg/L	Ethylbenzene µg/L	Xylenes µg/L	Trimethybenzenes µg/L	MTBE µg/L
WDNR ES	5	1000	700	10000	480	60
9/15/98	<0.50	<0.60	<0.60	<1.7	<1.7	<0.92
12/2/98	<0.50	<0.60	<0.60	<1.7	<1.7	<0.92
3/30/99	<0.50	<0.60	<0.60	<1.7	<1.7	<0.92
6/10/99	<0.50	<0.60	<0.60	<1.7	<1.7	<0.92
9/20/99	<0.50	<0.60	<0.60	<1.7	<1.7	<0.92
12/3/99	<0.50	<0.60	<0.60	<1.7	<1.7	<0.92

= ES exceedance

WEISENBERGER TIE & LUMBER COMPANY
PVOC Analytical Results

MONITORING WELL		DPZ-3						
Date Sampled	Benzene (ug/L)	Toluene (ug/L)	Ethylbenzene (ug/L)	Xylenes (ug/L)	1,2,4 TMB (ug/L)	1,3,5 TMB (ug/L)	MTBE (ug/L)	comments
WDNR ES	5	343	700	620			60	
06/03/92	<1.0	<1.0	<1.0	<3.0	<1.0	<1.0	NA	
08/20/92	<1.0	2.3	<1.0	11	6.2	<1.0	NA	
12/20/94	<1.0	<1.0	<1.0	4.2	<1.0	<1.0	NA	
03/11/98	<0.13	<0.20	<0.22	<0.23	<0.22	<0.29	<0.16	
06/24/98	<0.13	<0.20	<0.22	<0.23	<0.22	<0.29	<0.16	

DPZ-3

Date Sampled	Benzene µg/L	Toluene µg/L	Ethylbenzene µg/L	Xylenes µg/L	Trimethylbenzenes µg/L	MTBE µg/L
WDNR ES	5	1000	700	10000	480	60
9/15/98	<0.50	<0.60	<0.60	<1.7	<1.7	<0.92
12/2/98	<0.50	<0.60	<0.60	<1.7	<1.7	<0.92
3/30/99	<0.50	<0.60	<0.60	<1.7	<1.7	<0.92
6/10/99	<0.50	<0.60	<0.60	<1.7	<1.7	<0.92
9/20/99	<0.50	<0.60	<0.60	<1.7	<1.7	<0.92
12/3/99	<0.50	<0.60	<0.60	1.8	<1.7	<0.92


= ES exceedance

WEISENBERGER TIE & LUMBER COMPANY
PVOC Analytical Results

MONITORING WELL		DPZ-4						
Date Sampled	Benzene (ug/L)	Toluene (ug/L)	Ethylbenzene (ug/L)	Xylenes (ug/L)	1,2,4 TMB (ug/L)	1,3,5 TMB (ug/L)	MTBE (ug/L)	comments
WDNR ES	5	343	700	620			60	
06/03/92	-	-	-	-	-	-	-	NS
08/20/92	-	-	-	-	-	-	-	NS
12/20/94	<1.0	2.4	<1.0	<3.0	<1.0	<1.0	NA	
03/11/98	<0.13	<0.20	<0.22	<0.23	<0.22	<0.29	<0.16	
06/24/98	<0.13	<0.20	<0.22	<0.23	<0.22	<0.29	<0.16	

DPZ-4

Date Sampled	Benzene µg/L	Toluene µg/L	Ethylbenzene µg/L	Xylenes µg/L	Trimethylbenzenes µg/L	MTBE µg/L
WDNR ES	5	1000	700	10000	480	60
9/15/98	<0.50	<0.60	<0.60	<1.7	<1.7	<0.92
12/2/98	<0.50	<0.60	<0.60	<1.7	<1.7	<0.92
3/30/99	<0.50	<0.60	<0.60	<1.7	<1.7	<0.92
6/10/99	<0.50	<0.60	<0.60	<1.7	<1.7	<0.92
9/20/99	<0.50	<0.60	<0.60	<1.7	<1.7	<0.92
12/3/99	<0.50	<0.60	<0.60	<1.7	<1.7	<0.92

 = ES exceedance

WEISENBERGER TIE & LUMBER COMPANY
PVOC Analytical Results

MONITORING WELL		DPZ-5						
Date Sampled	Benzene (ug/L)	Toluene (ug/L)	Ethylbenzene (ug/L)	Xylenes (ug/L)	1,2,4 TMB (ug/L)	1,3,5 TMB (ug/L)	MTBE (ug/L)	comments
WDNR ES	5	343	700	620			60	
06/03/92	-	-	-	-	-	-	-	NS
08/20/92	<1.0	<1.0	<1.0	<3.0	<1.0	<1.0	NA	
12/20/94	<1.0	<1.0	<1.0	<3.0	<1.0	<1.0	NA	
03/11/98	-	-	-	-	-	-	-	NS
06/24/98	-	-	-	-	-	-	-	NS

DPZ-5

Date Sampled	Benzene µg/L	Toluene µg/L	Ethylbenzene µg/L	Xylenes µg/L	Trimethybenzenes µg/L	MTBE µg/L
WDNR ES	5	1000	700	10000	480	60
3/6/00	<0.50	<0.60	<0.60	<1.7	<1.7	<0.92

= ES exceedance

WEISENBERGER TIE & LUMBER COMPANY
PVOC Analytical Results

MONITORING WELL		DPZ-6						
Date Sampled	Benzene (ug/L)	Toluene (ug/L)	Ethylbenzene (ug/L)	Xylenes (ug/L)	1,2,4 TMB (ug/L)	1,3,5 TMB (ug/L)	MTBE (ug/L)	comments
WDNR ES	5	343	700	620			60	
06/03/92	-	-	-	-	-	-	-	NS
08/20/92	-	-	-	-	-	-	-	NS
12/20/94	<1.0	<1.0	<1.0	<3.0	<1.0	<1.0	NA	
03/11/98	<0.13	<0.20	<0.22	<0.23	<0.22	<0.29	<0.16	
06/24/98	<0.13	<0.20	<0.22	<0.23	<0.22	<0.29	<0.16	

DPZ-6

Date Sampled	Benzene µg/L	Toluene µg/L	Ethylbenzene µg/L	Xylenes µg/L	Trimethybenzenes µg/L	MTBE µg/L
WDNR ES	5	1000	700	10000	480	60
9/15/98	<0.50	0.83	<0.60	<1.7	<1.7	<0.92
12/2/98	Not Sampled					
3/30/99	<0.50	<0.60	<0.60	<1.7	<1.7	<0.92
6/10/99	<0.50	<0.60	<0.60	<1.7	<1.7	<0.92
9/20/99	<0.50	<0.60	<0.60	<1.7	<1.7	<0.92
12/3/99	<0.50	<0.60	<0.60	<1.7	<1.7	<0.92

= ES exceedance

PVOC							
3/11/98	Benzene	Toluene	Ethylbenzene	Xylene	1,2,4 TMB	1,3,5 TMB	MTBE
WDNR ES	5	343	700	620			60
DPW-1	<0.13	<0.20	<0.22	<0.23	<0.22	<0.29	<0.16
DPW-2	<0.13	<0.20	<0.22	3.7	4.1	1.0	0.79
DPW-3	0.31	<0.20	<0.22	6.1	2.0	0.66	0.76
DPW-4	0.16	<0.20	<0.22	0.63	<0.22	<0.29	1.5
DPW-5	<0.13	<0.20	<0.22	<0.23	<0.22	<0.29	<0.16
DPW-6	<0.13	<0.20	<0.22	<0.23	<0.22	<0.29	<0.16
DPW-7	<0.13	<0.20	<0.22	<0.23	<0.22	<0.29	<0.16
DPW-8	<0.13	<0.20	<0.22	<0.23	<0.22	<0.29	0.37

PVOC							
6/23/98	Benzene	Toluene	Ethylbenzene	Xylene	1,2,4 TMB	1,3,5 TMB	MTBE
WDNR ES	5	343	700	620			60
DPW-1	<0.13	<0.20	<0.22	<0.23	<0.22	<0.29	<0.16
DPW-2	<0.13	<0.20	<0.22	3.3	6.4	1.4	<0.16
DPW-3	0.66	1.4	2.0	24	21	3.2	<0.16
DPW-4	<0.13	<0.20	<0.22	<0.23	<0.22	<0.29	<0.16
DPW-5	<0.13	<0.20	<0.22	<0.23	<0.22	<0.29	<0.16
DPW-6	<0.13	<0.20	<0.22	<0.23	<0.22	<0.29	<0.16
DPW-7	<0.13	<0.20	<0.22	<0.23	<0.22	<0.29	<0.16
DPW-8	<0.13	<0.20	<0.22	<0.23	<0.22	<0.29	<0.16

NOTES:

ug/L = micrograms per liter

- = no analytical

1,2,4 TMB = 1,2,4 Trimethylbenzene

1,3,5 TMB = 1,3,5 Trimethylbenzene

MTBE = Methyl-tert-butyl ether

= indicates exceedance to WDNR enforcement standards (ES)

MONITORING WELLS
SVOC ANALYSIS

WEISENBERGER TIE & LUMBER COMPANY SVOC Analytical Results

MONITORING WELL MW-2										
Date Sampled (mm/dd/yy)	Naphthalene (ug/L)	Acenaphthylene (ug/L)	Acenaphthene (ug/L)	Fluorene (ug/L)	Penta. (ug/L)	2-Methylphenol (ug/L)	Phenanthrene (ug/L)	Fluoranthene (ug/L)	Pyrene (ug/L)	Comments
	40				1.6					
06/03/92	<10	<10	<10	<10	<50	<10	<10	<10	<10	
08/20/92	<12	<12	<12	<12	<58	<12	<12	<12	<12	
12/20/94	<10	<10	<10	<10	<25	<10	<10	<10	<10	
03/14/95	<11	<11	<11	<11	<26	<11	<11	<11	<11	
06/20/95	<10	<10	<10	<10	<25	<10	<10	<10	<10	
09/12/95	<11	<11	<11	<11	<53	<11	<11	<11	<11	
12/13/95	<10	<10	<10	<10	<10	<10	<25	<10	<10	
03/06/96	<10	<10	<10	<10	<25	<10	<10	<10	<10	
06/12/96	<10	<10	<10	<10	<25	<10	<10	<10	<10	
09/18/96	<10	<10	<10	<10	<25	<10	<10	<10	<10	
12/17/96	<10	<10	<10	<10	<25	<10	<10	<10	<10	
03/18/97	<2.0	<1.4	<1.3	<1.5	<2.3	<0.94	<0.69	<0.87	<0.66	
12/17/97	-	-	-	-	-	-	-	-	-	NS
03/11/98	-	-	-	-	-	-	-	-	-	NS
06/23/98										

= ES exceedance

WEISENBERGER TIE & LUMBER COMPANY SVOC Analytical Results

MONITORING WELL		MW-3								
Date Sampled (mm/dd/yy)	Naphthalene (ug/L)	Acenaphthylene (ug/L)	Acenaphthene (ug/L)	Fluorene (ug/L)	Penta. (ug/L)	2-Methylphenol (ug/L)	Phenanthrene (ug/L)	Fluoranthene (ug/L)	Pyrene (ug/L)	Comments
WDNR ES	40				1.0					
06/03/92	330	<110	<110	<110	37000	<110	<110	<110	<110	
08/20/92	<1000	<1000	<1000	<1000	31000	<1000	<1000	<1000	<1000	
12/20/94	170	<1000	<1000	<1000	22000	<1000	<1000	<1000	<1000	
03/15/95	300	<1000	<1000	<1000	32000	<1000	<1000	<1000	<1000	
06/21/95	<1000	<1000	<1000	<1000	28000	<1000	<1000	<1000	<1000	
09/14/95	180	5	6	12	24000	<10	<10	<10	<10	
12/14/95	290	<2000	<2000	<2000	34000	<2000	<2000	<2000	<2000	
03/06/96	370	<2000	<2000	<2000	34000	<2000	<2000	<2000	<2000	
06/13/96	260	<500	<500	<500	20000	<500	<500	<500	<500	
Dup (6/13/96)	250	<500	<500	<500	19000	<500	<500	<500	<500	
09/19/96	<2000	<2000	<2000	<2000	19000	<2000	<2000	<2000	<2000	
Dup (9/19/96)	<2000	<2000	<2000	<2000	19000	<2000	<2000	<2000	<2000	
12/17/96	-	-	-	-	-	-	-	-	-	
03/18/97	<400	<280	<260	<300	23000	<190	<140	<170	<130	
09/10/97	68	<7.0	<6.5	<7.5	18000	<4.7	<3.4	<4.3	<3.3	
Dup(9/10/97)	49	<7.0	<6.5	<7.5	18000	<4.7	<3.4	<4.3	<3.3	
12/17/97	<210	<140	<130	<150	15000	<97	<71	<90	<68	
03/11/98	260	<210	<230	<240	12100	<270	<240	<240	<240	
06/23/98	220	<10	16	17	7400	<10	17	<10	<10	

MW-3

Date Sampled	Naphthalene µg/L	Acenaphthylene µg/L	Acenaphthene µg/L	Fluorene µg/L	Pentachlorophenol µg/L	2-Methylphenol µg/L	Phenanthrene µg/L	Fluoranthene µg/L	Pyrene µg/L
WDNR ES	40			400	1.0			400	250
9/15/98	100	<0.66	3.7	<0.82	6840	<0.94	5.5	<0.96	<0.42
12/2/98	173	<0.66	2.8	7.7	12900	<0.94	13	<0.96	<0.42
3/30/99	113	<6.6	<8.4	<8.2	10600	<9.4	7.4	<9.6	<4.2
6/10/99	63	<6.6	<8.4	<8.2	9760	<9.4	7.4	<9.6	<4.2
9/20/99	129	<6.6	<8.4	<8.2	13000	<9.4	<6.8	<9.6	<4.2
12/3/99	169	<6.6	<8.4	8.4	13300	<9.4	10	<9.6	<4.2
3/6/00	146	<15	<11	<12	18600	<17	<14	<20	<22
6/30/00	34	<15	<11	<12	13900	<17	<14	<20	<22
9/27/00	163	<15	<11	<12	19600	<17	<14	<20	<22
12/27/00	151	<2.9	7.8	8.2	23700	<3.4	10	<4.0	<4.4
3/28/01	<14	<15	<11	<12	14900	<17	<14	<20	<22
6/27/01	60	<1.5	7.3	11	5830	<1.7	<1.4	<2.0	<2.2

= ES exceedance

WEISENBERGER TIE & LUMBER COMPANY SVOC Analytical Results

MONITORING WELL MW-5										
Date Sampled (mm/dd/yy)	Naphthalene (ug/L)	Acenaphthylene (ug/L)	Acenaphthene (ug/L)	Fluorene (ug/L)	Penta. (ug/L)	2-Methylphenol (ug/L)	Phenanthrene (ug/L)	Fluoranthene (ug/L)	Pyrene (ug/L)	Comments
WDNR ES	40				1.0					
06/03/92	170	<10	<10	<10	9900	<10	7	<10	<10	
08/20/92	160	<40	<40	<40	11000	<40	<40	<40	<40	
12/20/94	370	<1000	<1000	<1000	24000	<1000	<1000	<1000	<1000	
03/15/95	160	<1000	<1000	<1000	11000	<1000	<1000	<1000	<1000	
06/21/95	<1000	<1000	<1000	<1000	12000	<1000	<1000	<1000	<1000	
09/13/95	160	<10	<10	<10	7800	<10	9	<10	<10	
12/14/95	<1000	<1000	<1000	<1000	11000	<1000	<1000	<1000	<1000	
DUP(12/14/95)	<1000	<1000	<1000	<1000	11000	<1000	<1000	<1000	<1000	
3/6/96	<1000	<1000	<1000	<1000	9100	<1000	<1000	<1000	<1000	
6/13/96	<500	<500	<500	<500	7700	<500	<500	<500	<500	
9/18/96	<500	<500	<500	<500	5600	<500	<500	<500	<500	
12/17/96	<10	<10	<10	<10	5000	<10	<10	<10	<10	
3/19/97	<200	<140	<130	<150	6700	<94	<69	<87	<66	
9/10/97	<2.0	<1.4	<1.3	<1.5	2.4	<0.94	<0.69	<1.5	<0.66	
12/17/97	<2.1	<1.4	<1.3	<1.5	74	<0.91	<0.71	<0.90	<0.68	
3/11/98	4.1	<2.1	<2.3	<2.4	1400	<2.6	<2.4	>2.4	<2.4	
06/23/98	<24	<21	<23	<24	1900	<26	<24	<24	<24	

MW-5

Date Sampled	Naphthalene µg/L	Acenaphthylene µg/L	Acenaphthene µg/L	Fluorene µg/L	Pentachlorophenol µg/L	2-Methylphenol µg/L	Phenanthrene µg/L	Fluoranthene µg/L	Pyrene µg/L
WDNR ES	40			400	1.0			400	250
9/15/98	45	<0.66	<0.84	1.1	3700	<0.94	3.4	<0.96	<0.42
12/2/98	72	<0.66	<0.84	1.3	4270	<0.94	4.9	<0.96	<0.42
3/30/99	60	<3.3	<4.2	<4.1	3190	<4.7	<3.4	<4.8	<2.1
6/10/99	<3.5	<3.3	<4.2	<4.1	2910	<4.7	<3.4	<4.8	<2.1
9/20/99	<3.5	<3.3	<4.2	<4.1	3860	<4.7	<3.4	<4.8	<2.1
12/3/99	53	<6.6	<8.4	<8.2	3470	<9.4	<6.8	<9.6	<4.2
3/6/00	29	<7.3	<5.6	<6.0	3530	<8.5	<7.1	<9.9	<11
6/30/00	<14	<15	<11	<12	3400	<17	<14	<20	<22
9/27/00	<14	<15	<11	<12	3150	<17	<14	<20	<22
12/27/00	7.2	<2.9	<2.2	<2.4	803	<3.4	<2.8	<4.0	<4.4
3/28/01	36	<15	<11	<12	4240	<17	<14	<20	<22
6/27/01	31	<1.5	<1.1	<1.2	2650	<1.7	2.5	<2.0	<2.2

= ES exceedance

WEISENBERGER TIE & LUMBER COMPANY SVOC Analytical Results

MONITORING WELL MW-6										
Date Sampled (mm/dd/yy)	Naphthalene (ug/L)	Acenaphthylene (ug/L)	Acenaphthene (ug/L)	Fluorene (ug/L)	Penta. (ug/L)	2-Methylphenol (ug/L)	Phenanthrene (ug/L)	Fluoranthene (ug/L)	Pyrene (ug/L)	Comments
WDNR ES	40				400	1.0			400	250
06/03/92	<11	<11	<11	<11	<54	<11	<11	<11	<11	
08/20/92	<10	<10	<10	<10	<50	<10	<10	<10	<10	
12/20/94	<10	<10	<10	<10	<25	<10	<10	<10	<10	
03/15/95	<10	<10	<10	<10	16	<10	<10	<10	<10	
06/21/95	<11	<11	<11	<11	23	<11	<11	<11	<11	
Dup (6/21/95)	<10	<10	<10	<10	32	<10	<10	<10	<10	
09/13/95	<10	<10	<10	<10	<50	<10	<10	<10	<10	
12/13/95	<10	<10	<10	<10	<10	<25	<10	<10	<10	
03/06/96	<10	<10	<10	<10	<10	<25	<10	<10	<10	
06/13/96	-	-	-	-	-	-	-	-	-	
09/19/96	<10	<10	<10	<10	<25	<10	<10	<10	<10	
12/17/96	-	-	-	-	-	-	-	-	-	NS
03/18/97	-	-	-	-	-	-	-	-	-	NS
12/17/97	-	-	-	-	-	-	-	-	-	NS
03/11/98	NS	NS	NS	NS	NS	NS	NS	NS	NS	Bent Casing
06/23/98	<2.4	<2.1	<2.3	<2.4	<3.0	<2.6	<2.4	<2.4	<2.4	

MW-6

Date Sampled	Naphthalene µg/L	Acenaphthylene µg/L	Acenaphthene µg/L	Fluorene µg/L	Pentachlorophenol µg/L	2-Methylphenol µg/L	Phenanthrene µg/L	Fluoranthene µg/L	Pyrene µg/L
WDNR ES	40			400	1.0			400	250
9/15/98	Not Sampled								
12/2/98	Not Sampled								
3/30/99	15	<0.66	<0.84	<0.82	475	<0.94	<0.68	<0.96	<0.42
6/10/99	<0.70	<0.66	<0.84	<0.82	<0.90	<0.94	<0.68	<0.96	<0.42
9/20/99	<0.70	<0.66	<0.84	<0.82	6.2	<0.94	<0.68	<0.96	<0.42
12/3/99	<0.70	<0.66	<0.84	<0.82	79	<0.94	<6.8	<0.96	<0.42
3/6/00	<1.4	<1.5	<1.1	<1.2	487	<1.7	<1.4	<2.0	<2.2
6/30/00	<1.4	<1.5	<1.1	<1.2	<0.9	<1.7	<1.4	<2.0	<2.2
9/27/00	<1.4	<1.5	<1.1	<1.2	2.4	<1.7	<1.4	<2.0	<2.2
12/27/00	<1.4	<1.5	<1.1	<1.2	28	<1.7	<1.4	<2.0	<2.2
3/28/01	2.3	<1.5	<1.1	<1.2	421	<1.7	<1.4	<2.0	<2.2
6/27/01	<1.4	<1.5	<1.1	<1.2	85	<1.7	<1.4	<2.0	<2.2

= ES exceedance

WEISENBERGER TIE & LUMBER COMPANY SVOC Analytical Results

MONITORING WELL MW-7										
Date Sampled (mm/dd/yy)	Naphthalene (ug/L)	Acenaphthylene (ug/L)	Acenaphthene (ug/L)	Fluorene (ug/L)	Penta. (ug/L)	2-Methylphenol (ug/L)	Phenanthrene (ug/L)	Fluoranthene (ug/L)	Pyrene (ug/L)	Comments
WDNR ES	40				1.0					
06/03/92	<11	<11	14	16	2900	<11	14	<11	<11	
08/20/92	18	<20	9	10	3000	<20	10	<20	<20	
12/20/94	29	<100	<100	<100	1300	<100	<100	<100	<100	
03/14/95	10	<100	<100	<100	2900	<100	10	<100	<100	
06/20/95	<10	<10	<10	<10	2300	<10	<10	<10	<10	
09/12/95	<10	<10	3	3	2800	<10	<10	<10	<10	
12/14/95	<100	<100	11	10	2800	<100	<100	<100	<100	
03/06/96	3	<10	1	2	360	<10	2	<10	<10	
06/13/96	<250	<250	<250	<250	2200	<250	<250	<250	<250	
09/18/96	<200	<200	<200	<200	2400	<200	<200	<200	<200	
12/17/96	72	<10	10	9.4	1800	<10	5	<10	<10	
03/19/97	<100	<70	<65	<75	2400	<47	<34	<44	<33	
09/10/97	<2.0	<1.4	7.5	<0.87	2300	<0.94	<0.69	<0.87	<0.66	
12/17/97	-	-	-	-	-	-	-	-	-	DRY
03/11/98	-	-	-	-	-	-	-	-	-	DRY
06/23/98	<2.4	<2.1	<2.3	<2.4	550	<2.6	<2.4	<2.4	<2.4	

MW-7

Date Sampled	Naphthalene µg/L	Acenaphthylene µg/L	Acenaphthene µg/L	Fluorene µg/L	Pentachlorophenol µg/L	2-Methylphenol µg/L	Phenanthrene µg/L	Fluoranthene µg/L	Pyrene µg/L
WDNR ES	40			400	1.0			400	250
9/15/98	39	<0.66	3.5	3.1	1120	<0.94	<0.68	<0.96	<0.42
12/2/98	20	<0.66	4.9	4.4	1210	<0.94	1.8	<0.96	<0.42
3/30/99	<0.70	<0.66	<0.84	<0.82	91	<0.94	<0.68	<0.96	<0.42
6/10/99	5.9	<0.66	4.1	3.5	795	<0.94	<0.68	<0.96	<0.42
9/20/99	53	<1.3	10	10	1360	<1.9	6.1	<1.9	<0.84
12/3/99	<3.5	<3.3	9.9	11	1380	<4.7	7.5	<4.8	<2.1
3/6/00	<6.9	<7.3	<5.6	<6.0	2090	<8.5	<7.1	<9.9	<11
6/30/00	<6.9	<7.3	<5.6	<6.0	818	<8.5	<7.1	<9.9	<11
9/27/00	46	<7.3	7.2	6.8	1320	<8.5	<7.1	<9.9	<11
12/27/00	51	<2.9	11	11	1830	<3.4	8.6	<4.0	<4.4
3/28/01	Not Sampled								
6/27/01	60	<1.5	7.7	6.3	1190	<1.7	1.6	<2.0	<2.2

= ES exceedance

WEISENBERGER TIE & LUMBER COMPANY SVOC Analytical Results

MONITORING WELL MW-10										Comments
Date Sampled (mm/dd/yy)	Naphthalene (ug/L)	Acenaphthylene (ug/L)	Acenaphthene	Fluorene (ug/L)	Penta. (ug/L)	2-Methylphenol (ug/L)	Phenanthrene (ug/L)	Fluoranthene (ug/L)	Pyrene (ug/L)	
WDNR ES	40				1.0					
06/03/92	25	<10	12	12	1500	<10	10	<10	<10	
08/20/92	<40	<40	<40	<40	730	<40	<40	<40	<40	
12/20/94	19	<20	7	8	436	<20	10	<20	<20	
Dup (12/20/94)	<10	<10	<10	<10	460	<20	10	<20	<20	
03/15/95	34	<20	9	10	1100	<20	11	<20	<20	
06/21/95	<11	<11	<11	<11	920	<11	<11	<11	<11	
Dup (6/21/95)	<10	<10	<10	<10	1100	<10	<10	<10	<10	
09/13/95	<10	<10	8	8	910	<10	5	<10	<10	
12/14/95	19	<20	12	<20	390	<20	18	<20	2	
3/6/96	NS	NS	NS	NS	NS	NS	NS	NS	NS	
6/13/96	<10	<10	2	2	100	<10	<10	<10	<10	
9/18/96	<10	<10	<10	<10	81	<10	<10	<10	<10	
12/17/96	<50	<50	6.7	8.9	150	<50	<50	<50	<50	
Dup (12/17/96)	10	<10	7.9	8.8	140	<10	9.3	<10	1.2	
3/19/97	<20	<14	<13	<15	400	<9.4	<6.9	<8.7	<6.6	
9/10/97	<20	<14	<13	<15	250	<9.4	<6.9	<8.7	<6.6	
12/17/97	<10	<7.2	<6.7	<7.7	180	<4.8	<3.6	<4.5	<3.4	
3/11/98	-	-	-	-	-	-	-	-	-	DRY
06/23/98	<2.4	<2.1	6.3	3.4	330	<2.6	<2.4	<2.4	<2.4	

MW-10

Date Sampled	Naphthalene µg/L	Acenaphthylene µg/L	Acenaphthene µg/L	Fluorene µg/L	Pentachlorophenol µg/L	2-Methylphenol µg/L	Phenanthrene µg/L	Fluoranthene µg/L	Pyrene µg/L
WDNR ES	40			400	1.0			400	250
9/15/98	2.6	<0.66	2.4	3.4	176	<0.94	3.1	<0.96	<0.42
12/2/98	8.2	<0.66	5.0	5.5	482	<0.94	5.3	<0.96	<0.42
3/30/99	7.2	<1.3	7.5	7.4	563	<1.9	5.3	<1.9	<0.8
6/10/99	2.6	<1.3	<1.7	2.1	221	<1.9	<1.4	<1.9	<0.8
9/20/99	<1.4	<1.3	<1.7	2.7	81	<1.9	<1.4	<1.9	<0.84
12/3/99	4.3	<1.3	4.0	4.4	153	<1.9	2.9	<1.9	<0.84
3/6/00	8.0	<1.5	2.9	2.8	832	<1.7	<1.4	<2.0	<3.0
6/30/00	2.5	<1.5	1.7	1.9	225	<1.7	<1.4	<2.0	<2.2
9/27/00	3.6	<1.5	2.5	3.2	266	<1.7	2.6	<2.0	<2.2
12/27/00	13	<2.9	11	13	550	<3.4	9.6	<4.0	<4.4
3/28/01	Not Sampled								
6/27/01	<1.4	<1.5	<1.1	<1.2	58	<1.7	<1.4	<2.0	<2.2

= ES exceedance

WEISENBERGER TIE & LUMBER COMPANY SVOC Analytical Results

MONITORING WELL DMW-1										
Date Sampled (mm/dd/yy)	Naphthalene (ug/L)	Acenaphthylene (ug/L)	Acenaphthene (ug/L)	Fluorene (ug/L)	Penta. (ug/L)	2-Methylphenol (ug/L)	Phenanthrene (ug/L)	Fluoranthene (ug/L)	Pyrene (ug/L)	Comments
WDNR ES	40				1.0					
06/03/92	<100	<100	<100	<100	17000	<100	<100	<100	<100	
Dup (6/03/92)	<100	<100	<100	<100	17000	<100	<100	<100	<100	
08/20/92	<500	<500	<500	<500	16000	<500	<500	<500	<500	
12/20/94	-	-	-	-	-	-	-	-	-	NS
03/15/95	11	<10	25	<10	6300	<10	75	6	9	
06/21/95	<10	<10	15	<10	3700	<10	49	5	8	
09/13/95	12	<10	21	31	12000	<10	40	6	<10	
12/14/95	<100	<100	15	<100	2800	<100	32	<100	<100	
Dup (12/14/95)	<200	<200	27	<200	4500	<200	60	<200	<200	
03/06/96	-	-	-	-	-	-	-	-	-	NS
6/13/96	<1000	<1000	<1000	<1000	14000	<1000	50	<1000	<1000	
9/19/96	<2000	<2000	<2000	<2000	12000	<2000	<2000	<2000	<2000	
Dup (9/19/96)	<2000	<2000	<2000	<2000	11000	<2000	<2000	<2000	<2000	
12/17/96	-	-	-	-	-	-	-	-	-	NS
3/18/97	-	-	-	-	-	-	-	-	-	NS
9/10/97	<100	<70	<65	<75	2400	<47	<34	<44	<33	
12/17/97	<100	<72	<67	<77	10000	<48	<36	<45	<34	
3/11/98	<240	<210	<230	<240	12300	<260	<240	<240	<240	
06/23/98	<10	<10	34	<10	11500	<10	<10	<10	<10	

DMW-1

Date Sampled	Naphthalene µg/L	Acenaphthylene µg/L	Acenaphthene µg/L	Fluorene µg/L	Pentachlorophenol µg/L	2-Methylphenol µg/L	Phenanthrene µg/L	Fluoranthene µg/L	Pyrene µg/L
WDNR ES	40			400	1.0			400	250
9/15/98	24	<0.66	16	26	7940	<0.94	40	3.6	6.7
12/2/98	28	<0.66	31	27	11200	<0.94	63	8.3	13
3/30/99	<7.0	<6.6	<8.4	24	6980	<9.4	51	<9.6	12
6/10/99	<7.0	<6.6	11	<8.2	3530	<9.4	12	<9.6	<4.2
9/20/99	<7.0	<6.6	16	15	6170	<9.4	25	<9.6	<4.2
12/3/99	14	<6.6	94	96	9590	<9.4	230	21	38
3/6/00	<6.9	<7.3	25	12	10300	<8.5	26	<9.9	<11
6/30/00	<14	<15	18	16	6530	<17	31	<20	<22
9/27/00	37	<15	57	72	10500	<17	134	<20	23
12/27/00	Not sampled								
3/28/01	<14	<15	26	23	11200	<17	39	<20	<22
6/27/01	<1.4	<1.5	24	24	4050	<1.7	29	4.2	5.4

= ES exceedance

WEISENBERGER TIE & LUMBER COMPANY
SVOC Analytical Results

MONITORING WELL DMW-2										
Date Sampled (mm/dd/yy)	Naphthalene (ug/L)	Acenaphthylene (ug/L)	Acenaphthene (ug/L)	Fluorene (ug/L)	Penta. (ug/L)	2-Methylphenol (ug/L)	Phenanthrene (ug/L)	Fluoranthene (ug/L)	Pyrene (ug/L)	Comments
WDNR ES	40				1.0					
3/11/98	-	-	-	-	-	-	-	-	-	
06/23/98	<2.4	<2.1	<2.3	<2.4	<3.0	<2.6	<2.4	<2.4	<2.4	

DMW-2


Date Sampled	Naphthalene µg/L	Acenaphthylene µg/L	Acenaphthene µg/L	Fluorene µg/L	Pentachlorophenol µg/L	2-Methylphenol µg/L	Phenanthrene µg/L	Fluoranthene µg/L	Pyrene µg/L
WDNR ES	40			400	1.0			400	250
9/15/98					Not Sampled				
12/2/98					Not Sampled				
3/30/99	<0.70	<0.66	<0.84	<0.82	1.7	<0.94	<0.68	<0.96	<0.42
6/10/99	<0.70	<0.66	<0.84	<0.82	2.2	<0.94	<0.68	<0.96	<0.42
9/20/99					Not Sampled				
12/3/99					Not Sampled				
6/10/99	<0.70	<0.66	<0.84	<0.82	2.2	<0.94	<0.68	<0.96	<0.42
3/6/00	<1.4	<1.5	<1.1	<1.2	<0.9	<1.7	<1.4	<2.0	<2.2
6/30/00	<1.4	<1.5	<1.1	<1.2	16	<1.7	<1.4	<2.0	<2.2
9/27/00	<1.4	<1.5	<1.1	<1.2	3.7	<1.7	<1.4	<2.0	<2.2
12/27/00					Not sampled				
3/28/01	<1.4	<1.5	<1.1	<1.2	<0.9	<1.7	<1.4	<2.0	<2.2
6/27/01	<1.4	<1.5	<1.1	<1.2	<1.7	<1.7	<1.4	<2.0	<2.2

= ES exceedance

WEISENBERGER TIE & LUMBER COMPANY
SVOC Analytical Results

DMW-3

Date Sampled	Naphthalene µg/L	Acenaphthylene µg/L	Acenaphthene µg/L	Fluorene µg/L	Pentachlorophenol µg/L	2-Methylphenol µg/L	Phenanthrene µg/L	Fluoranthene µg/L	Pyrene µg/L
WDNR ES	40			400	1.0			400	250
3/6/00									Not Sampled
6/30/00									Not Sampled
9/27/00									Not Sampled
12/27/00									Not Sampled
3/28/01									Not Sampled
6/27/01									Not Sampled

 = ES exceedance

WEISENBERGER TIE & LUMBER COMPANY SVOC Analytical Results

MONITORING WELL DMW-4										
Date Sampled (mm/dd/yy)	Naphthalene (ug/L)	Acenaphthylene (ug/L)	Acenaphthene (ug/L)	Fluorene (ug/L)	Penta. (ug/L)	2-Methylphenol (ug/L)	Phenanthrene (ug/L)	Fluoranthene (ug/L)	Pyrene (ug/L)	Comments
WDNR ES	40				1.4					
06/03/92	<10	<10	10	10	7100	<10	7	<10	<10	
08/20/92	<20	<20	<20	<20	5700	<20	<20	<20	<20	
12/20/94	-	-	-	-	-	-	-	-	-	NS
03/14/95	-	-	-	-	-	-	-	-	-	NS
06/21/95	<10	<10	<10	<10	<25	<10	<10	<10	<10	
09/13/95	<10	<10	<10	<10	130	<10	<10	<10	<10	
12/13/95	-	-	-	-	-	-	-	-	-	NS
03/06/96	-	-	-	-	-	-	-	-	-	NS
06/13/96	<10	<10	<10	<10	1	<10	<10	<10	<10	
09/18/96	-	-	-	-	-	-	-	-	-	NS
12/17/96	-	-	-	-	-	-	-	-	-	NS
03/18/97	-	-	-	-	-	-	-	-	-	MUD
12/17/97	-	-	-	-	-	-	-	-	-	DRY
03/11/98	-	-	-	-	-	-	-	-	-	DRY
06/23/98	<2.4	<2.1	<2.3	<2.4	<3.0	<2.6	<2.4	<2.4	<2.4	

DMW-4

Date Sampled	Naphthalene µg/L	Acenaphthylene µg/L	Acenaphthene µg/L	Fluorene µg/L	Pentachlorophenol µg/L	2-Methylphenol µg/L	Phenanthrene µg/L	Fluoranthene µg/L	Pyrene µg/L
WDNR ES	40			400	1.0			400	250
9/15/98	<0.70	<0.66	<0.84	<0.82	16	<0.94	<0.68	<0.96	<0.42
12/2/98	Not Sampled								
3/30/99	<0.70	<0.66	<0.84	<0.82	3.7	<0.94	<0.68	<0.96	<0.42
6/10/99	<0.70	<0.66	<0.84	<0.82	12	<0.94	<0.68	<0.96	<0.42
9/20/99	3.7	<0.66	<0.84	1.3	2050	<0.94	<0.68	<0.96	<0.42
12/3/99	Not Sampled								
3/6/00	Not Sampled								
6/30/00	<1.4	<1.5	<1.1	<1.2	3.0	<1.7	<1.4	<2.0	<2.2
9/27/00	<1.4	<1.5	<1.1	<1.2	43	<1.7	<1.4	<2.0	<2.2
12/27/00	Not Sampled								
3/28/01	<1.4	<1.5	<1.1	<1.2	4.2	<1.7	<1.4	<2.0	<2.2
6/27/01	<1.4	<1.5	<1.1	<1.2	38	<1.7	<1.4	<2.0	<2.2

= ES exceedance

WEISENBERGER TIE & LUMBER COMPANY
SVOC Analytical Results

MONITORING WELL		DMW-5								Comments
Date Sampled (mm/dd/yy)	Naphthalene (ug/L)	Acenaphthylene (ug/L)	Acenaphthene (ug/L)	Fluorene (ug/L)	Penta. (ug/L)	2-Methylphenol (ug/L)	Phenanthrene (ug/L)	Fluoranthene (ug/L)	Pyrene (ug/L)	
WDNR ES	40					1.0				
08/03/92	<11	<11	<11	<11	<57	<11	<11	<11	<11	
08/20/92	<10	<10	<10	<10	<50	<10	<10	<10	<10	
12/20/94	<10	<10	<10	<10	<25	<10	<10	<10	<10	
03/14/95	-	-	-	-	-	-	-	-	-	NS
06/22/95	<10	<10	<10	<10	<25	<10	<10	<10	<10	
09/13/95	<10	<10	<10	<10	<50	<10	<10	<10	<10	
12/13/95	-	-	-	-	-	-	-	-	-	NS
03/06/96	-	-	-	-	-	-	-	-	-	NS
06/12/96	-	-	-	-	-	-	-	-	-	NS
09/18/96	-	-	-	-	-	-	-	-	-	NS
12/17/96	-	-	-	-	-	-	-	-	-	NS
03/18/97	-	-	-	-	-	-	-	-	-	NS
12/17/97	-	-	-	-	-	-	-	-	-	DRY
03/11/98	-	-	-	-	-	-	-	-	-	DRY
06/23/98	<2.4	<2.1	<2.3	<2.4	<3.0	<2.6	<2.4	<2.4	<2.4	

DMW-5

Date Sampled	Naphthalene µg/L	Acenaphthylene µg/L	Acenaphthene µg/L	Fluorene µg/L	Pentachlorophenol µg/L	2-Methylphenol µg/L	Phenanthrene µg/L	Fluoranthene µg/L	Pyrene µg/L
WDNR ES	40			400	1.0			400	250
9/15/98	<0.70	<0.66	<0.84	<0.82	<0.90	<0.94	<0.68	<0.96	<0.42
12/2/98	<0.70	<0.66	<0.84	<0.82	39	<0.94	<0.68	<0.96	<0.42
3/30/99	<0.70	<0.66	<0.84	<0.82	1.0	<0.94	<0.68	<0.96	<0.42
6/10/99	<0.70	<0.66	<0.84	<0.82	<0.90	<0.94	<0.68	<0.96	<0.42
9/20/99	<0.70	<0.66	<0.84	<0.82	<0.90	<0.94	<0.68	<0.96	<0.42
12/3/99	<0.70	<0.66	<0.84	<0.82	18	<0.94	<0.68	<0.96	<0.42
3/6/00	Not Sampled								
6/30/00	<1.4	<1.5	<1.1	<1.2	0.98	<1.7	<1.4	<2.0	<2.2
9/27/00	<1.4	<1.5	<1.1	<1.2	1.5	<1.7	<1.4	<2.0	<2.2
12/27/00	Not Sampled								
3/28/01	Not Sampled								
6/27/01	<1.4	<1.5	<1.1	<1.2	<1.7	<1.7	<1.4	<2.0	<2.2

 = ES exceedance

WEISENBERGER TIE & LUMBER COMPANY SVOC Analytical Results

MONITORING WELL DMW-6a										
Date Sampled (mm/dd/yy)	Naphthalene (ug/L)	Acenaphthylene (ug/L)	Acenaphthene (ug/L)	Fluorene (ug/L)	Penta. (ug/L)	2-Methylphenol (ug/L)	Phenanthrene (ug/L)	Fluoranthene (ug/L)	Pyrene (ug/L)	Comments
WDNR ES	40				400	1.0			400	250
06/03/92	<11	<11	<11	<11	600	<11	<11	<11	<11	
08/20/92	<10	<10	<10	<10	110	<10	<10	<10	<10	
12/20/94	3	<20	<20	<20	330	<20	<20	<20	<20	
Dup (12/20/94)	<20	<20	<20	<20	370	<20	<20	<20	<20	
03/14/95	-	-	-	-	-	-	-	-	-	NS
06/20/95	<11	<11	<11	<11	38	<11	<11	<11	<11	
09/12/95	<10	<10	<10	<10	<50	<10	<10	<10	<10	
12/13/95	-	-	-	-	-	-	-	-	-	NS
03/05/96	-	-	-	-	-	-	-	-	-	NS
06/12/96	-	-	-	-	-	-	-	-	-	NS
09/18/96	-	-	-	-	-	-	-	-	-	NS
12/17/96	-	-	-	-	-	-	-	-	-	NS
03/18/97	-	-	-	-	-	-	-	-	-	DRY
12/17/97	-	-	-	-	-	-	-	-	-	DRY
03/11/98	-	-	-	-	-	-	-	-	-	DRY
06/23/98	<2.4	<2.1	<2.3	<2.4	<3.0	<2.6	<2.4	<2.4	<2.4	

DMW-6A

Date Sampled	Naphthalene µg/L	Acenaphthylene µg/L	Acenaphthene µg/L	Fluorene µg/L	Pentachlorophenol µg/L	2-Methylphenol µg/L	Phenanthrene µg/L	Fluoranthene µg/L	Pyrene µg/L
WDNR ES	40			400	1.0			400	250
9/15/98	<0.70	<0.66	<0.84	<0.82	3.6	<0.94	<0.68	<0.96	<0.42
12/2/98	<0.70	<0.66	<0.84	<0.82	6.5	<0.94	<0.68	<0.96	<0.42
3/30/99	Not Sampled								
6/10/99	<0.70	<0.66	<0.84	<0.82	<0.90	<0.94	<0.68	<0.96	<0.42
9/20/99	<0.70	<0.66	<0.84	<0.82	2.1	<0.94	<0.68	<0.96	<0.42
12/3/99	Not Sampled								
3/6/00	Not Sampled								
6/30/00	<1.4	<1.5	<1.1	<1.2	2.5	<1.7	<1.4	<2.0	<2.2
9/27/00	<1.4	<1.5	<1.1	<1.2	2.4	<1.7	<1.4	<2.0	<2.2
12/27/00	Not Sampled								
3/28/01	Not Sampled								
6/27/01	<1.4	<1.5	<1.1	<1.2	<1.7	<1.7	<1.4	<2.0	<2.2

= ES exceedance

WEISENBERGER TIE & LUMBER COMPANY SVOC Analytical Results

MONITORING WELL DMW-7										
Date Sampled (mm/dd/yy)	Naphthalene (ug/L)	Acenaphthylene (ug/L)	Acenaphthene (ug/L)	Fluorene (ug/L)	Penta. (ug/L)	2-Methylphenol (ug/L)	Phenanthrene (ug/L)	Fluoranthene (ug/L)	Pyrene (ug/L)	Comments
WDNR ES	40				1.0					
06/03/92	-	-	-	-	-	-	-	-	-	NS
08/20/92	<10	<10	<10	<10	<50	<10	<10	<10	<10	
12/20/94	<50	<50	<10	<50	1100	<50	<50	<50	<50	
03/14/95	<50	<50	<50	<50	1500	<50	<50	<50	<50	
06/20/95	<10	<10	<10	<10	590	<10	<10	<10	<10	
09/13/95	<10	<10	<10	<10	23	<10	<10	<10	<10	
12/13/95	<10	<10	<10	<10	53	<10	<10	<10	<10	
03/06/96	<10	<10	<10	<10	8	<10	<10	<10	<10	
06/12/96	<10	<10	<10	<10	25	<10	<10	<10	<10	
09/18/96	<10	<10	<10	<10	25	<10	<10	<10	<10	
12/17/96	-	-	-	-	-	-	-	-	-	NS
03/18/97	-	-	-	-	-	-	-	-	-	NS
12/17/97	-	-	-	-	-	-	-	-	-	NS
03/11/98	<2.4	<2.1	<2.3	<2.4	<3.0	<2.6	<2.4	<2.4	<2.4	
06/23/98	<2.4	<2.1	<2.3	<2.4	<3.0	<2.6	<2.4	<2.4	<2.4	

DMW-7

Date Sampled	Naphthalene µg/L	Acenaphthylene µg/L	Acenaphthene µg/L	Fluorene µg/L	Pentachlorophenol µg/L	2-Methylphenol µg/L	Phenanthrene µg/L	Fluoranthene µg/L	Pyrene µg/L
WDNR ES	40			400	1.0			400	250
9/15/98	<0.70	<0.66	<0.84	<0.82	<0.90	<0.94	<0.68	<0.96	<0.42
12/2/98	<0.70	<0.66	<0.84	<0.82	1.0	<0.94	<0.68	<0.96	<0.42
3/30/99	<0.70	<0.66	<0.84	<0.82	1.3	<0.94	<0.68	<0.96	<0.42
6/10/99	<0.70	<0.66	<0.84	<0.82	<0.90	<0.94	<0.68	<0.96	<0.42
9/20/99	<0.70	<0.66	<0.84	<0.82	2.0	<0.94	<0.68	<0.96	<0.42
12/3/99	<0.70	<0.66	<0.84	<0.82	0.90	<0.94	<0.68	<0.96	<0.42
3/6/00	<1.4	<1.5	<1.1	<1.2	<0.9	<1.7	<1.4	<2.0	<2.2
6/30/00	<1.4	<1.5	<1.1	<1.2	<0.9	<1.7	<1.4	<2.0	<2.2
9/27/00	<1.4	<1.5	<1.1	<1.2	<0.9	<1.7	<1.4	<2.0	<2.2
12/27/00	<1.4	<1.5	<1.1	<1.2	<0.9	<1.7	<1.4	<2.0	<2.2
3/28/01	<1.4	<1.5	<1.1	<1.2	<0.9	<1.7	<1.4	<2.0	<2.2
6/27/01	<1.4	<1.5	<1.1	<1.2	<1.7	<1.7	<1.4	<2.0	<2.2


= ES exceedance

WEISENBERGER TIE & LUMBER COMPANY
SVOC Analytical Results

MONITORING WELL DMW-8										
Date Sampled (mm/dd/yy)	Naphthalene (ug/L)	Acenaphthylene (ug/L)	Acenaphthene (ug/L)	Fluorene (ug/L)	Penta. (ug/L)	2-Methylphenol (ug/L)	Phenanthrene (ug/L)	Fluoranthene (ug/L)	Pyrene (ug/L)	Comments
WDNR ES	40				1.0					
06/03/92	-	-	-	-	-	-	-	-	-	NS
08/20/92	<10	<10	<10	<10	<50	<10	<10	<10	<10	
12/20/94	<10	<10	<10	<10	<25	<10	<10	<10	<10	
03/14/95	<10	<10	<10	<10	<25	<10	<10	<10	<10	
06/20/95	<10	<10	<10	<10	<25	<10	<10	<10	<10	
09/13/95	<11	<11	<11	<11	<53	<11	<11	<11	<11	
12/14/95	<10	<10	<10	<10	<25	<10	<10	<10	<10	
03/06/96	<10	<10	<10	<10	<25	<10	<10	<10	<10	
06/12/96	<10	<10	<10	<10	<25	<10	<10	<10	<10	
09/18/96	<10	<10	<10	<10	<25	<10	<10	<10	<10	
12/17/96	-	-	-	-	-	-	-	-	-	NS
03/18/97	-	-	-	-	-	-	-	-	-	NS
12/17/97	-	-	-	-	-	-	-	-	-	NS
03/11/98	<2.4	<2.1	<2.3	<2.4	<3.0	<2.6	<2.4	<2.4	<2.4	
06/23/98	<2.4	<2.1	<2.3	<2.4	<3.0	<2.6	<2.4	<2.4	<2.4	

DMW-8

Date Sampled	Naphthalene µg/L	Acenaphthylene µg/L	Acenaphthene µg/L	Fluorene µg/L	Pentachlorophenol µg/L	2-Methylphenol µg/L	Phenanthrene µg/L	Fluoranthene µg/L	Pyrene µg/L
WDNR ES	40			400	1.0			400	250
9/15/98	<0.70	<0.66	<0.84	<0.82	<0.90	<0.94	<0.68	<0.96	<0.42
12/2/98	<0.70	<0.66	<0.84	<0.82	<0.90	<0.94	<0.68	<0.96	<0.42
3/30/99	<0.70	<0.66	<0.84	<0.82	<0.90	<0.94	<0.68	<0.96	<0.42
6/10/99	<0.70	<0.66	<0.84	<0.82	<0.90	<0.94	<0.68	<0.96	<0.42
9/20/99	<0.70	<0.66	<0.84	<0.82	<0.90	<0.94	<0.68	<0.96	<0.42
12/3/99	<0.70	<0.66	<0.84	<0.82	<0.90	<0.94	<0.68	<0.96	<0.42

 = ES exceedance

WEISENBERGER TIE & LUMBER COMPANY
SVOC Analytical Results

DMW-10

Date Sampled	Naphthalene µg/L	Acenaphthylene µg/L	Acenaphthene µg/L	Fluorene µg/L	Pentachlorophenol µg/L	2-Methylphenol µg/L	Phenanthrene µg/L	Fluoranthene µg/L	Pyrene µg/L
WDNR ES	40			400	1.0			400	250
3/6/00	<1.4	<1.5	<1.1	<1.2	<0.90	<1.7	<1.4	<2.0	<2.2
6/30/00	<1.4	<1.5	<1.1	<1.2	<0.9	<1.7	<1.4	<2.0	<2.2
9/27/00	<1.4	<1.5	<1.1	<1.2	<0.9	<1.7	<1.4	<2.0	<2.2
12/27/00	<1.4	<1.5	<1.1	<1.2	0.98	<1.7	<1.4	<2.0	<2.2
3/28/01	<1.4	<1.5	<1.1	<1.2	<0.90	<1.7	<1.4	<2.0	<2.2
6/27/01	<1.4	<1.5	<1.1	<1.2	<1.7	<1.7	<1.4	<2.0	<2.2


= ES exceedance

WEISENBERGER TIE & LUMBER COMPANY
SVOC Analytical Results

MONITORING WELL DMW-12										
Date Sampled (mm/dd/yy)	Naphthalene (ug/L)	Acenaphthylene (ug/L)	Acenaphthene (ug/L)	Fluorene (ug/L)	Penta. (ug/L)	2-Methylphenol (ug/L)	Phenanthrene (ug/L)	Fluoranthene (ug/L)	Pyrene (ug/L)	Comments
WDNR ES	40				1.0					
06/03/92	-	-	-	-	-	-	-	-	-	NS
08/20/92	-	-	-	-	-	-	-	-	-	NS
12/20/94	<10	<10	<10	<10	<25	<10	<10	<10	<10	
03/14/95	<10	<10	<10	<10	<25	<10	<10	<10	<10	
06/20/95	<10	<10	<10	<10	<25	<10	<10	<10	<10	
09/12/95	<10	<10	<10	<10	<52	<10	<10	<10	<10	
12/13/95	-	-	-	-	-	-	-	-	-	DRY
03/06/96	-	-	-	-	-	-	-	-	-	NS
06/13/96	-	-	-	-	-	-	-	-	-	DRY
09/18/96	-	-	-	-	-	-	-	-	-	DRY
12/17/96	-	-	-	-	-	-	-	-	-	DRY
03/18/97	-	-	-	-	-	-	-	-	-	DRY
12/17/97	-	-	-	-	-	-	-	-	-	DRY
03/11/98	-	-	-	-	-	-	-	-	-	DRY
06/23/98	<2.4	<2.1	<2.3	<2.4	<3.0	<2.6	<2.4	<2.4	<2.4	

DMW-12

Date Sampled	Naphthalene µg/L	Acenaphthylene µg/L	Acenaphthene µg/L	Fluorene µg/L	Pentachlorophenol µg/L	2-Methylphenol µg/L	Phenanthrene µg/L	Fluoranthene µg/L	Pyrene µg/L
WDNR ES	40			400	1.0			400	250
9/15/98	<0.70	<0.66	<0.84	<0.82	<0.90	<0.94	<0.68	<0.96	<0.42
12/2/98	<0.70	<0.66	<0.84	<0.82	<0.90	<0.94	<0.68	<0.96	<0.42
3/30/99	<0.70	<0.66	<0.84	<0.82	<0.90	<0.94	<0.68	<0.96	<0.42
6/10/99	<0.70	<0.66	<0.84	<0.82	<0.90	<0.94	<0.68	<0.96	<0.42
9/20/99	<0.70	<0.66	<0.84	<0.82	<0.90	<0.94	<0.68	<0.96	<0.42
12/3/99	<0.70	<0.66	<0.84	<0.82	<0.90	<0.94	<0.68	<0.96	<0.42

 = ES exceedance

WEISENBERGER TIE & LUMBER COMPANY
SVOC Analytical Results

MONITORING WELL DMW-13										
Date Sampled (mm/dd/yy)	Naphthalene (ug/L)	Acenaphthylene (ug/L)	Acenaphthene (ug/L)	Fluorene (ug/L)	Penta. (ug/L)	2-Methylphenol (ug/L)	Phenanthrene (ug/L)	Fluoranthene (ug/L)	Pyrene (ug/L)	Comments
WDNR ES	40				1.0					
09/12/95	<10	<10	<10	<10	<50	<10	<10	<10	<10	
12/13/95	<10	<10	<10	<10	<25	<10	<10	<10	<10	
03/06/96	-	-	-	-	-	-	-	-	-	FROZEN
06/12/96	<10	<10	<10	<10	<25	<10	<10	<10	<10	
09/18/96	<10	<10	<10	<10	<25	<10	<10	<10	<10	
12/17/96	-	-	-	-	-	-	-	-	-	NS
03/18/97	-	-	-	-	-	-	-	-	-	FROZEN
12/17/97	-	-	-	-	-	-	-	-	-	NS
03/11/98	<2.4	<2.1	<2.3	<2.4	<3.0	<2.6	<2.4	<2.4	<2.4	
06/23/98	<2.4	<2.1	<2.3	<2.4	<3.0	<2.6	<2.4	<2.4	<2.4	

DMW-13

Date Sampled	Naphthalene µg/L	Acenaphthylene µg/L	Acenaphthene µg/L	Fluorene µg/L	Pentachlorophenol µg/L	2-Methylphenol µg/L	Phenanthrene µg/L	Fluoranthene µg/L	Pyrene µg/L
WDNR ES	40			400	1.0			400	250
9/15/98	<0.70	<0.66	<0.84	<0.82	<0.90	<0.94	<0.68	<0.96	<0.42
12/2/98	<0.70	<0.66	<0.84	<0.82	22	<0.94	<0.68	<0.96	<0.42
3/30/99	<0.70	<0.66	<0.84	<0.82	<0.90	<0.94	<0.68	<0.96	<0.42
6/10/99	<0.70	<0.66	<0.84	<0.82	<0.90	<0.94	<0.68	<0.96	<0.42
9/20/99	<0.70	<0.66	<0.84	<0.82	1.7	<0.94	<0.68	<0.96	<0.42
12/3/99	<0.70	<0.66	<0.84	<0.82	<0.90	<0.94	<0.68	<0.96	<0.42
3/6/00	<1.4	<1.5	<1.1	<1.2	<0.90	<1.7	<1.4	<2.0	<2.2
6/30/00	<1.4	<1.5	<1.1	<1.2	<0.9	<1.7	<1.4	<2.0	<2.2
9/27/00	<1.4	<1.5	<1.1	<1.2	<0.9	<1.7	<1.4	<2.0	<2.2
12/27/00	<1.4	<1.5	<1.1	<1.2	<0.9	<1.7	<1.4	<2.0	<2.2
3/28/01	Not Sampled								
6/27/01	<1.4	<1.5	<1.1	<1.2	<1.7	<1.7	<1.4	<2.0	<2.2

= ES exceedance

WEISENBERGER TIE & LUMBER COMPANY SVOC Analytical Results

PIEZOMETER DPZ-1										
Date Sampled (mm/dd/yy)	Naphthalene (ug/L)	Acenaphthylene (ug/L)	Acenaphthene (ug/L)	Fluorene (ug/L)	Penta. (ug/L)	2-Methylphenol (ug/L)	Phenanthrene (ug/L)	Fluoranthene (ug/L)	Pyrene (ug/L)	Comments
WDNR ES	40				1.0					
06/03/92	<40	<40	91	120	12000	<40	260	19	24	
08/20/92	<200	<200	<200	<200	5600	<200	<200	<200	<200	
12/20/94	-	-	-	-	-	-	-	-	-	NS
03/15/95	16	<54	19	27	7900	<54	39	<54	6	
06/22/95	<50	<50	27	31	5500	<50	39	<50	17	
09/14/95	<10	<10	8	5	5100	<10	<10	2	4	
12/14/95	<250	<250	<250	<250	5700	<250	29	<250	<250	
03/06/96	<250	<250	28	<250	9000	<250	33	<250	13	
06/13/96	<1000	<1000	<1000	<1000	5700	<1000	<1000	<1000	<1000	
Dup (6/13/96)	<1000	<1000	<1000	<1000	5300	<1000	<1000	<1000	<1000	
09/19/96	<1000	<1000	<1000	<1000	5600	<1000	<1000	<1000	<1000	
12/17/96	<1000	<1000	<1000	<1000	6700	<1000	<1000	<1000	<1000	
03/19/97	<200	<140	<130	<150	4900	<94	<69	<87	<66	
09/10/97	<20	<14	<13	<15	5000	<9.4	<6.9	<15	<6.6	
12/17/97	<200	<140	<130	<150	3900	<94	<69	<87	<66	
03/11/98	<51	<45	<49	<51	1300	<55	<51	<51	<51	
06/23/98	<2.4	<2.1	<2.3	<2.4	2300	<2.6	<2.4	<2.4	<2.4	

DPZ-1

Date Sampled	Naphthalene µg/L	Acenaphthylene µg/L	Acenaphthene µg/L	Fluorene µg/L	Pentachlorophenol µg/L	2-Methylphenol µg/L	Phenanthrene µg/L	Fluoranthene µg/L	Pyrene µg/L
WDNR ES	40			400	1.0			400	250
9/15/98	101	<0.66	<0.84	<0.82	6480	<0.94	<0.68	<0.96	<0.42
12/2/98	129	<0.66	<0.84	2.9	7500	<0.94	<0.68	<0.96	<0.42
3/30/99	59	<3.3	<4.2	<4.1	4460	<4.7	<3.4	<4.8	<2.1
6/10/99	<7.0	<0.66	<0.84	<8.2	3960	<0.94	<0.68	<0.96	<0.42
9/20/99	<7.0	<6.6	<8.4	<8.2	5830	<9.4	<6.8	<9.6	<4.2
12/3/99	96	<6.6	<8.4	<8.2	4450	<9.4	<6.8	<9.6	<4.2
3/6/00	191	<1.5	<1.1	1.5	8300	<1.7	2.1	<2.0	<2.2
6/30/00	<14	<15	<11	<12	6910	<17	<14	<20	<22
9/27/00	113	<15	<11	<12	7000	<17	<14	<20	<22
12/27/00	112	<2.9	<2.2	<2.4	11000	<3.4	<2.8	<4.0	<4.4
3/28/01	39	<15	<11	<12	7990	<17	<14	<20	<22
6/27/01	2.18	<1.5	<1.1	<1.2	3120	<1.7	<1.4	<2.0	<2.2


= ES exceedance

WEISENBERGER TIE & LUMBER COMPANY
SVOC Analytical Results

PIEZOMETER		DPZ-1a								
Date Sampled (mm/dd/yy)	Naphthalene (ug/L)	Acenaphthylene (ug/L)	Acenaphthene (ug/L)	Fluorene (ug/L)	Penta. (ug/L)	2-Methylphenol (ug/L)	Phenanthrene (ug/L)	Fluoranthene (ug/L)	Pyrene (ug/L)	Comments
WDNR ES	40				1.0					
06/03/92	-	-	-	-	-	-	-	-	-	NS
08/20/92	<10	<10	<10	<10	130	<10	<10	<10	<10	
12/20/94	<10	<10	<10	<10	<25	<10	<10	<10	<10	
03/14/95	-	-	-	-	-	-	-	-	-	NS
06/22/95	<10	<10	<10	<10	<25	<10	<10	<10	<10	
09/13/95	-	-	-	-	-	-	-	-	-	NS
12/13/95	-	-	-	-	-	-	-	-	-	NS
03/05/96	-	-	-	-	-	-	-	-	-	NS
06/13/96	-	-	-	-	-	-	-	-	-	NS
09/19/96	<11	<11	<11	<11	4.0	<11	<11	<11	<11	
12/17/96	-	-	-	-	-	-	-	-	-	NS
03/18/97	-	-	-	-	-	-	-	-	-	NS
12/17/97	-	-	-	-	-	-	-	-	-	NS
03/11/98	<2.4	<2.1	<2.3	<2.4	<3.0	<2.6	<2.4	<2.4	<2.4	
06/23/98	<2.4	<2.1	<2.3	<2.4	<3.0	<2.6	<2.4	<2.4	<2.4	

DPZ-1a

Date Sampled	Naphthalene µg/L	Acenaphthylene µg/L	Acenaphthene µg/L	Fluorene µg/L	Pentachlorophenol µg/L	2-Methylphenol µg/L	Phenanthrene µg/L	Fluoranthene µg/L	Pyrene µg/L
WDNR ES	40			400	1.0			400	250
9/15/98	<0.70	<0.66	<0.84	<0.82	3.6	<0.94	<0.68	<0.96	<0.42
12/2/98	<0.70	<0.66	<0.84	<0.82	12	<0.94	<0.68	<0.96	<0.42
3/30/99	<0.70	<0.66	<0.84	<0.82	2.2	<0.94	<0.68	<0.96	<0.42
6/10/99	<0.70	<0.66	<0.84	<0.82	4.0	<0.94	<0.68	<0.96	<0.42
9/20/99	<0.70	<0.66	<0.84	<0.82	8.6	<0.94	<0.68	<0.96	<0.42
12/3/99	<0.70	<0.66	<0.84	<0.82	11	<0.94	<0.68	<0.96	<0.42
3/6/00	<1.4	<1.5	<1.1	<1.2	6.4	<1.7	<1.4	<2.0	<2.2
6/30/00	<1.4	<1.5	<1.1	<1.2	4.1	<1.7	<1.4	<2.0	<2.2
9/27/00	<1.4	<1.5	<1.1	<1.2	4.9	<1.7	<1.4	<2.0	<2.2
12/27/00	<1.4	<1.5	<1.1	<1.2	6.2	<1.7	<1.4	<2.0	<2.2
3/28/01	<1.4	<1.5	<1.1	<1.2	8.1	<1.7	<1.4	<2.0	<2.2
6/27/01	<1.4	<1.5	<1.1	<2.0	<1.7	<1.7	<1.4	<2.0	<2.2


 = ES exceedance

WEISENBERGER TIE & LUMBER COMPANY
SVOC Analytical Results

PIEZOMETER DPZ-2										
Date Sampled (mm/dd/yy)	Naphthalene (ug/L)	Acenaphthylene (ug/L)	Acenaphthene (ug/L)	Fluorene (ug/L)	Penta. (ug/L)	2-Methylphenol (ug/L)	Phenanthrene (ug/L)	Fluoranthene (ug/L)	Pyrene (ug/L)	Comments
WDNR ES	40				1.0					
06/03/92	<10	<10	<10	<10	<53	<10	<10	<10	<10	
Dup (6/03/92)	<10	<10	<10	<10	<53	<10	<10	<10	<10	
08/20/92	<10	<10	<10	<10	<50	<10	<10	<10	<10	
Dup (8/20/92)	<10	<10	<10	<10	<50	<10	<10	<10	<10	
12/20/94	<10	<10	<10	<10	<25	<10	<10	<10	<10	
03/15/95	<10	<10	<10	<10	<25	<10	<10	<10	<10	
06/22/95	<10	<10	<10	<10	<25	<10	<10	<10	<10	
09/13/95	<10	<10	<10	<10	<51	<10	<10	<10	<10	
12/13/95	<10	<10	<10	<10	<25	<10	<10	<10	<10	
03/06/96	<10	<10	<10	<10	<25	<10	<10	<10	<10	
06/13/96	<10	<10	<10	<10	<25	<10	<10	<10	<10	
09/19/96	<10	<10	<10	<10	<25	<10	<10	<10	<10	
12/17/96	-	-	-	-	-	-	-	-	-	NS
03/18/97	-	-	-	-	-	-	-	-	-	NS
12/17/97	-	-	-	-	-	-	-	-	-	NS
03/11/98	<2.4	<2.1	<2.3	<2.4	<3.0	<2.6	<2.4	<2.4	<2.4	
06/23/98	<2.4	<2.1	<2.3	<2.4	<3.0	<2.6	<2.4	<2.4	<2.4	

DPZ-2

Date Sampled	Naphthalene µg/L	Acenaphthylene µg/L	Acenaphthene µg/L	Fluorene µg/L	Pentachlorophenol µg/L	2-Methylphenol µg/L	Phenanthrene µg/L	Fluoranthene µg/L	Pyrene µg/L
WDNR ES	40			400	1.0			400	250
9/15/98	<0.70	<0.66	<0.84	<0.82	11	<0.94	<0.68	<0.96	<0.42
12/2/98	<0.70	<0.66	<0.84	<0.82	9.5	<0.94	<0.68	<0.96	<0.42
3/30/99	<0.70	<0.66	<0.84	<0.82	8.3	<0.94	<0.68	<0.96	<0.42
6/10/99	<0.70	<0.66	<0.84	<0.82	<0.90	<0.94	<0.68	<0.96	<0.42
9/20/99	<0.70	<0.66	<0.84	<0.82	4.8	<0.94	<0.68	<0.96	<0.42
12/3/99	<0.70	<0.66	<0.84	<0.82	28	<0.94	<0.68	<0.96	<0.42
3/6/00	1.8	<1.5	<1.1	<1.2	666	<1.7	<1.4	<2.0	<2.2
6/30/00	<1.4	<1.5	<1.1	<1.2	<0.9	<1.7	<1.4	<2.0	<2.2
9/27/00	<1.4	<1.5	<1.1	<1.2	4.9	<1.7	<1.4	<2.0	<2.2
12/27/00	<1.4	<1.5	<1.1	<1.2	1.8	<1.7	<1.4	<2.0	<2.2
3/28/01	3.8	<1.5	<1.1	<1.2	984	<1.7	<1.4	<2.0	<2.2
6/27/01	<1.4	<1.5	<1.1	<1.2	1.9	<1.7	<1.4	<2.0	<2.2

 = ES exceedance

WEISENBERGER TIE & LUMBER COMPANY SVOC Analytical Results

PIEZOMETER		DPZ-3								
Date Sampled (mm/dd/yy)	Naphthalene (ug/L)	Acenaphthylene (ug/L)	Acenaphthene (ug/L)	Fluorene (ug/L)	Penta. (ug/L)	2-Methylphenol (ug/L)	Phenanthrene (ug/L)	Fluoranthene (ug/L)	Pyrene (ug/L)	Comments
WDNR ES	40				1.4					
06/03/92	<11	<11	<11	<11	2000	<11	<11	<11	<11	
08/20/92	<10	<10	<10	<10	2100	<10	<10	<10	<10	
12/20/94	<100	<100	<100	<100	1500	<100	<100	<100	<100	
Dup (12/20/94)	<100	<100	<100	<100	1500	<100	3	<20	<100	
03/14/95	<100	<100	<100	<100	1800	<100	<100	<100	<100	
Dup (3/14/95)	8	<20	<20	<20	1600	<20	<10	<10	<20	
06/20/95	<11	<11	<11	<11	1500	<11	<11	<11	<11	
Dup (6/20/95)	<10	<10	<10	<10	1400	<10	<10	<10	<10	
09/12/95	8	<10	<10	<10	1200	<10	2	<100	<10	
12/14/95	<100	<100	<100	<100	840	<100	<100	<20	<100	
03/06/96	<20	<20	<20	<20	210	<20	<20	<10	<20	
06/13/96	<10	<10	<10	<10	<25	<10	<100	<100	<10	
09/18/96	<10	<10	<10	<10	<25	<10	<10	<10	<10	
12/17/96	-	-	-	-	-	-	-	-	-	NS
03/18/97	-	-	-	-	-	-	-	-	-	NS
9/10/97	<2.1	<1.5	<1.4	<1.6	15	<0.93	<0.73	<0.92	<0.69	
12/17/97	<2.1	<1.5	<1.4	<1.6	2.4	<0.98	<0.72	<0.91	<0.69	
3/11/98	<2.5	<2.2	<2.2	<2.5	<3.2	<2.7	<2.5	<2.5	<2.5	
06/23/98	<2.4	<2.1	<2.3	<2.4	<3.0	<2.6	<2.4	<2.4	<2.4	

DPZ-3

Date Sampled	Naphthalene µg/L	Acenaphthylene µg/L	Acenaphthene µg/L	Fluorene µg/L	Pentachlorophenol µg/L	2-Methylphenol µg/L	Phenanthrene µg/L	Fluoranthene µg/L	Pyrene µg/L
WDNR ES	40			400	1.0			400	250
9/15/98	<0.70	<0.66	<0.84	<0.82	<0.90	<0.94	<0.68	<0.96	<0.42
12/2/98	<0.70	<0.66	<0.84	<0.82	8.8	<0.94	<0.68	<0.96	<0.42
3/30/99	<0.70	<0.66	<0.84	<0.82	4.3	<0.94	<0.68	<0.96	<0.42
6/10/99	<0.70	<0.66	<0.84	<0.82	<0.90	<0.94	<0.68	<0.96	<0.42
9/20/99	<0.70	<0.66	<0.84	<0.82	36	<0.94	<0.68	<0.96	<0.42
12/3/99	<0.70	<0.66	<0.84	<0.82	369	<0.94	<0.68	<0.96	<0.42
3/6/00	<1.4	<1.5	<1.1	<1.2	318	<1.7	<1.4	<2.0	<2.2
6/30/00	<1.4	<1.5	<1.1	<1.2	1.4	<1.7	<1.4	<2.0	<2.2
9/27/00	<1.4	<1.5	<1.1	<1.2	54	<1.7	<1.4	<2.0	<2.2
12/27/00	<2.8	<2.9	<2.2	<2.4	72	<3.4	<2.8	<4.0	<4.4
3/28/01	>1.4	<1.5	<1.1	<1.2	257	<1.7	<1.4	<2.0	<2.2
6/27/01	<1.4	<1.5	<1.1	<1.2	<1.7	<1.7	<1.4	<2.0	<2.2

= ES exceedance

WEISENBERGER TIE & LUMBER COMPANY
SVOC Analytical Results

PIEZOMETER		DPZ-4								
Date Sampled (mm/dd/yy)	Naphthalene (ug/L)	Acenaphthylene (ug/L)	Acenaphthene (ug/L)	Fluorene (ug/L)	Penta. (ug/L)	2-Methylphenol (ug/L)	Phenanthrene (ug/L)	Fluoranthene (ug/L)	Pyrene (ug/L)	Comments
WDNR ES	40				400	1.0			400	250
06/03/92	-	-	-	-	-	-	-	-	-	NS
08/20/92	-	-	-	-	-	-	-	-	-	NS
12/20/94	<10	<10	<10	<10	<25	<10	<10	<10	<10	
03/15/95	<37	<37	<37	<37	47	<37	<37	<37	<37	
06/21/95	-	-	-	-	-	-	-	-	-	NS
09/13/95	-	<10	<10	<10	56	<10	<10	<10	<10	
12/13/95	-	<10	<10	<10	70	<10	<10	<10	<10	
03/06/96	-	-	-	-	-	-	-	-	-	NS
06/13/96	<10	<10	<10	<10	12	<10	<10	<10	<10	
09/18/96	<10	<10	<10	<10	12	<10	<10	<10	<10	
12/17/96	-	-	-	-	-	-	-	-	-	NS
03/18/97	-	-	-	-	-	-	-	-	-	NS
09/10/97	<2.0	<1.4	<1.3	<1.5	39	<0.94	<0.69	<0.87	<0.66	
12/17/97	-	-	-	-	-	-	-	-	-	DRY
03/11/98	<2.4	<2.1	<2.3	<2.4	13	<2.6	<2.4	<2.4	<2.4	
06/23/98	<2.4	<2.1	<2.3	<2.4	<3.0	<2.6	<2.4	<2.4	<2.4	

DPZ-4

Date Sampled	Naphthalene µg/L	Acenaphthylene µg/L	Acenaphthene µg/L	Fluorene µg/L	Pentachlorophenol µg/L	2-Methylphenol µg/L	Phenanthrene µg/L	Fluoranthene µg/L	Pyrene µg/L
WDNR ES	40			400	1.0			400	250
9/15/98	<0.70	<0.66	<0.84	<0.82	19	<0.94	<0.68	<0.96	<0.42
12/2/98	<0.70	<0.66	<0.84	<0.82	16	<0.94	<0.68	<0.96	<0.42
3/30/99	<0.70	<0.66	<0.84	<0.82	22	<0.94	<0.68	<0.96	<0.42
6/10/99	<0.70	<0.66	<0.84	<0.82	5.1	<0.94	<0.68	<0.96	<0.42
9/20/99	<0.70	<0.66	<0.84	<0.82	12	<0.94	<0.68	<0.96	<0.42
12/3/99	<0.70	<0.66	0.84	<0.82	9.9	<0.94	<0.68	<0.96	<0.42
3/6/00	<1.4	<1.5	<1.1	<1.2	27	<1.7	<1.4	<2.0	<2.2
6/30/00	<1.4	<1.5	<1.1	<1.2	6.9	<1.7	<1.4	<2.0	<2.2
9/27/00	<1.4	<1.5	<1.1	<1.2	7.6	<1.7	<1.4	<2.0	<2.2
12/27/00	Not Sampled								
3/28/01	<1.4	<1.5	<1.1	<1.2	<0.90	<1.7	<1.4	<2.0	<2.2
6/27/01	<1.4	<1.5	<1.1	<1.2	5.2	<1.7	<1.4	<2.0	<2.2


= ES exceedance

WEISENBERGER TIE & LUMBER COMPANY
SVOC Analytical Results

PIEZOMETER DPZ-5										
Date Sampled (mm/dd/yy)	Naphthalene (ug/L)	Acenaphthylene (ug/L)	Acenaphthene (ug/L)	Fluorene (ug/L)	Penta. (ug/L)	2-Methylphenol (ug/L)	Phenanthrene (ug/L)	Fluoranthene (ug/L)	Pyrene (ug/L)	Comments
WDNRES	40				1.0					
06/03/92	-	-	-	-	-	-	-	-	-	
08/20/92	<10	<10	<10	<10	<50	<10	<10	<10	<10	
12/20/94	<10	<10	<10	<10	220	<10	<10	<10	<10	
03/14/95	<10	<10	<10	<10	170	<10	<10	<10	<10	
Dup (3/14/95)	<20	<20	<20	<20	180	<20	<20	<20	<20	
06/21/95	<10	<10	<10	<10	<25	<10	<10	<10	<10	
09/13/95	<10	<10	<10	<10	180	<10	<10	<10	<10	
12/14/95	<10	<10	<10	<10	120	<10	<10	<10	<10	
3/6/96	<10	<10	<10	<10	120	<10	<10	<10	<10	
6/12/96	<10	<10	<10	<10	86	<10	<10	<10	<10	
9/18/96	<10	<10	<10	<10	70	<10	<10	<10	<10	
12/17/96	<10	<10	<10	<10	4.5	<10	<10	<10	<10	
3/18/97	<2.0	<1.4	<1.3	<1.5	<2.3	<0.94	<0.69	<0.87	<0.66	
9/10/97	<2.0	<1.4	<1.3	<1.5	<2.3	<0.94	<0.69	<0.87	<0.66	
12/17/97	-	-	-	-	-	-	-	-	-	NS
03/11/98	<2.4	<2.1	<2.3	<2.4	<3.0	<2.6	<2.4	<2.4	<2.4	
06/23/98	<2.4	<2.1	<2.3	<2.4	<3.0	<2.6	<2.4	<2.4	<2.4	

DPZ-5

Date Sampled	Naphthalene µg/L	Acenaphthylene µg/L	Acenaphthene µg/L	Fluorene µg/L	Pentachlorophenol µg/L	2-Methylphenol µg/L	Phenanthrene µg/L	Fluoranthene µg/L	Pyrene µg/L
WDNRES	40			400	1.0			400	250
9/15/98	<0.70	<0.66	<0.84	<0.82	<0.90	<0.94	<0.68	<0.96	<0.42
12/2/98	<0.70	<0.66	<0.84	<0.82	<0.90	<0.94	<0.68	<0.96	<0.42
3/30/99	<0.70	<0.66	<0.84	<0.82	<0.90	<0.94	<0.68	<0.96	<0.42
6/10/99	<0.70	<0.66	<0.84	<0.82	<0.90	<0.94	<0.68	<0.96	<0.42
9/20/99	<0.70	<0.66	<0.84	<0.82	<0.90	<0.94	<0.68	<0.96	<0.42
12/3/99	<0.70	<0.66	<0.84	<0.82	<0.90	<0.94	<0.68	<0.96	<0.42
3/6/00	<1.4	<1.5	<1.1	<1.2	<0.90	<1.7	<1.4	<2.0	<2.2
9/27/00	<1.4	<1.5	<1.1	<1.2	<0.90	<1.7	<1.4	<2.0	<2.2
12/27/00	<1.4	<1.5	<1.1	<1.2	<0.90	<1.7	<1.4	<2.0	<2.2
3/28/01	<1.4	<1.5	<1.1	<1.2	<0.90	<1.7	<1.4	<2.0	<2.2
6/27/01	<1.4	<1.5	<1.1	<1.2	<1.7	<1.7	<1.4	<2.0	<2.2

 = ES exceedance

WEISENBERGER TIE & LUMBER COMPANY
SVOC Analytical Results

PIEZOMETER		DPZ-6									Comments
Date Sampled (mm/dd/yy)	Naphthalene (ug/L)	Acenaphthylene (ug/L)	Acenaphthene (ug/L)	Fluorene (ug/L)	Penta. (ug/L)	2-Methylphenol (ug/L)	Phenanthrene (ug/L)	Fluoranthene (ug/L)	Pyrene (ug/L)		
WDR ES	40				1.0						
06/03/92	-	-	-	-	-	-	-	-	-	NS	
08/20/92	-	-	-	-	-	-	-	-	-	NS	
12/20/94	<20	<20	<20	<20	470	<20	<20	<20	<20		
03/14/95	<10	<10	<10	<10	<25	<10	<10	<10	<10		
06/20/95	<10	<10	<10	<10	<25	<10	<10	<10	<10		
09/13/95	<10	<10	<10	<10	<50	<10	<10	<10	<10		
12/13/95	-	-	-	-	-	-	-	-	-	NS	
03/06/96	<10	<10	<10	<10	<25	<10	<10	<10	<10		
06/12/96	-	-	-	-	-	-	-	-	-	NS	
09/18/96	-	-	-	-	-	-	-	-	-	NS	
12/17/96	-	-	-	-	-	-	-	-	-	NS	
03/18/97	-	-	-	-	-	-	-	-	-	NS	
12/17/97	-	-	-	-	-	-	-	-	-	NS	
03/11/98	<2.4	<2.1	<2.3	<2.4	<3.0	<2.6	<2.4	<2.4	<2.4		
06/23/98	<2.4	<2.1	<2.3	<2.4	<3.0	<2.6	<2.4	<2.4	<2.4		

DPZ-6

Date Sampled	Naphthalene µg/L	Acenaphthylene µg/L	Acenaphthene µg/L	Fluorene µg/L	Pentachlorophenol µg/L	2-Methylphenol µg/L	Phenanthrene µg/L	Fluoranthene µg/L	Pyrene µg/L
WDR ES	40			400	1.0			400	250
9/15/98	<0.70	<0.66	<0.84	<0.82	<0.90	<0.94	<0.68	<0.96	<0.42
12/2/98	Not Sampled								
3/30/99	<0.70	<0.66	<0.84	<0.82	<0.90	<0.94	<0.68	<0.96	<0.42
6/10/99	<0.70	<0.66	<0.84	<0.82	<0.90	<0.94	<0.68	<0.96	<0.42
9/20/99	<0.70	<0.66	<0.84	<0.82	<0.90	<0.94	<0.68	<0.96	<0.42
12/3/99	<0.70	<0.66	<0.84	<0.82	<0.90	<0.94	<0.68	<0.96	0.50
3/6/00	Not Sampled								
9/27/00	<1.4	<1.5	<1.1	<1.2	9.6	<1.7	<1.4	<2.0	<2.2
12/27/00	Not Sampled								
3/28/01	<1.4	<1.5	<1.1	<1.2	<0.90	<1.7	<1.4	<2.0	<2.2
6/27/01	<1.4	<1.5	<1.1	<1.2	<1.7	<1.7	<1.4	<2.0	<2.2

= ES exceedance

SEMI-VOLATILE GROUND WATER ANALYTICAL RESULTS

Weisenberger Tie and Lumber Company

Marathon City, Wisconsin

PUMPING WELL DPW-1										
Date Sampled (mm/dd/yy)	Naphthalene (ug/L)	Acenaphthylene (ug/L)	Acenaphthene (ug/L)	Fluorene (ug/L)	Penta. (ug/L)	2-Methylphenol (ug/L)	Phenanthrene (ug/L)	Fluoranthene (ug/L)	Pyrene (ug/L)	Comments
WDNR ES	40				1.9					
6/16/97	-	-	-	-	62	-	-	-	-	
12/17/97	-	-	-	-	43.0	-	-	-	-	
3/11/98	-	-	-	-	47.0	-	-	-	-	
06/23/98	<2.4	<2.1	<2.3	<2.4	6.8	<2.6	<2.4	<2.4	<2.4	

PUMPING WELL DPW-2										
Date Sampled (mm/dd/yy)	Naphthalene (ug/L)	Acenaphthylene (ug/L)	Acenaphthene (ug/L)	Fluorene (ug/L)	Penta. (ug/L)	2-Methylphenol (ug/L)	Phenanthrene (ug/L)	Fluoranthene (ug/L)	Pyrene (ug/L)	Comments
WDNR ES	40				1.9					
6/24/97	-	-	-	-	2900	-	-	-	-	
12/17/97	-	-	-	-	1200	-	-	-	-	
3/11/98	-	-	-	-	530	-	-	-	-	
06/23/98	12	<2.1	<2.3	4.5	910	<2.6	2.7	<2.4	<2.4	

PUMPING WELL DPW-3										
Date Sampled (mm/dd/yy)	Naphthalene (ug/L)	Acenaphthylene (ug/L)	Acenaphthene (ug/L)	Fluorene (ug/L)	Penta. (ug/L)	2-Methylphenol (ug/L)	Phenanthrene (ug/L)	Fluoranthene (ug/L)	Pyrene (ug/L)	Comments
WDNR ES	40				1.9					
6/16/97	-	-	-	-	3800	-	-	-	-	
12/17/97	-	-	-	-	3300	-	-	-	-	
3/11/98	-	-	-	-	2500	-	-	-	-	
06/23/98	<2.4	<2.1	2.9	2.6	4100	<2.6	<2.4	<2.4	<2.4	

PUMPING WELL DPW-4										
Date Sampled (mm/dd/yy)	Naphthalene (ug/L)	Acenaphthylene (ug/L)	Acenaphthene (ug/L)	Fluorene (ug/L)	Penta. (ug/L)	2-Methylphenol (ug/L)	Phenanthrene (ug/L)	Fluoranthene (ug/L)	Pyrene (ug/L)	Comments
WDNR ES	40				1.9					
6/16/97	-	-	-	-	3200	-	-	-	-	
12/17/97	-	-	-	-	2800	-	-	-	-	
3/11/98	-	-	-	-	3000	-	-	-	-	
06/23/98	<2.4	<2.1	<2.3	<2.4	270	<2.6	<2.4	<2.4	<2.4	

NOTES:

ug/L = micrograms per liter

- = no analysis

Penta. = Pentachlorophenol

Shaded value = NR 140 ES exceedance

NS = not sampled

SEMI-VOLATILE GROUND WATER ANALYTICAL RESULTS

Weisenberger Tie and Lumber Company

Marathon City, Wisconsin

PUMPING WELL DPW-5										
Date Sampled (mm/dd/yy)	Naphthalene (ug/L)	Acenaphthylene (ug/L)	Acenaphthene (ug/L)	Fluorene (ug/L)	Penta. (ug/L)	2-Methylphenol (ug/L)	Phenanthrene (ug/L)	Fluoranthene (ug/L)	Pyrene (ug/L)	Comments
WDNR ES	40	-	-	-	1.0	-	-	-	-	
6/16/97	-	-	-	-	<2.3	-	-	-	-	
12/17/97	-	-	-	-	-	-	-	-	-	NA
3/11/98	-	-	-	-	-	-	-	-	-	NA
06/23/98	-	-	-	-	-	-	-	-	-	NA

PUMPING WELL DPW-6										
Date Sampled (mm/dd/yy)	Naphthalene (ug/L)	Acenaphthylene (ug/L)	Acenaphthene (ug/L)	Fluorene (ug/L)	Penta. (ug/L)	2-Methylphenol (ug/L)	Phenanthrene (ug/L)	Fluoranthene (ug/L)	Pyrene (ug/L)	Comments
WDNR ES	40	-	-	-	1.0	-	-	-	-	
6/16/97	-	-	-	-	<2.3	-	-	-	-	
12/17/97	-	-	-	-	-	-	-	-	-	NA
3/11/98	-	-	-	-	-	-	-	-	-	NA
06/23/98	-	-	-	-	-	-	-	-	-	NA

PUMPING WELL DPW-7										
Date Sampled (mm/dd/yy)	Naphthalene (ug/L)	Acenaphthylene (ug/L)	Acenaphthene (ug/L)	Fluorene (ug/L)	Penta. (ug/L)	2-Methylphenol (ug/L)	Phenanthrene (ug/L)	Fluoranthene (ug/L)	Pyrene (ug/L)	Comments
WDNR ES	40	-	-	-	1.0	-	-	-	-	
6/16/97	-	-	-	-	<2.3	-	-	-	-	
12/17/97	-	-	-	-	-	-	-	-	-	NA
3/11/98	-	-	-	-	-	-	-	-	-	NA
06/23/98	-	-	-	-	-	-	-	-	-	NA

PUMPING WELL DPW-8										
Date Sampled (mm/dd/yy)	Naphthalene (ug/L)	Acenaphthylene (ug/L)	Acenaphthene (ug/L)	Fluorene (ug/L)	Penta. (ug/L)	2-Methylphenol (ug/L)	Phenanthrene (ug/L)	Fluoranthene (ug/L)	Pyrene (ug/L)	Comments
WDNR ES	40	-	-	-	1.0	-	-	-	-	
6/16/97	-	-	-	-	2.4	-	-	-	-	
12/17/97	-	-	-	-	<2.4	-	-	-	-	
3/11/98	-	-	-	-	-	-	-	-	-	NA
06/23/98	-	-	-	-	-	-	-	-	-	NA

NOTES:

ug/L = micrograms per liter

- = no analysis

Penta. = Pentachlorophenol

Shaded value = NR 140 ES exceedance

NS = not sampled

PRIVATE WELL ANALYSIS

WEISENBERGER TIE & LUMBER COMPANY
Private Well Analytical Results

Date Sampled	PCP ($\mu\text{g/L}$)
12/2/98	<0.04
3/30/99	0.3
6/10/99	<0.04
9/20/99	<0.04
12/3/99	<1.0
3/6/00	0.05
6/30/00	<0.04
9/27/00	0.07
12/27/00	0.12
3/28/01	0.20
6/27/01	0.05

 = Exceedance of WDNR Enforcement Standard of 1.0 $\mu\text{g/L}$

**CONTOUR MAPS
AND
FIELD DATA**

SITE NAME: WEISENBERGER TIE & LUMBER

DATE BAILED: 6/27/01

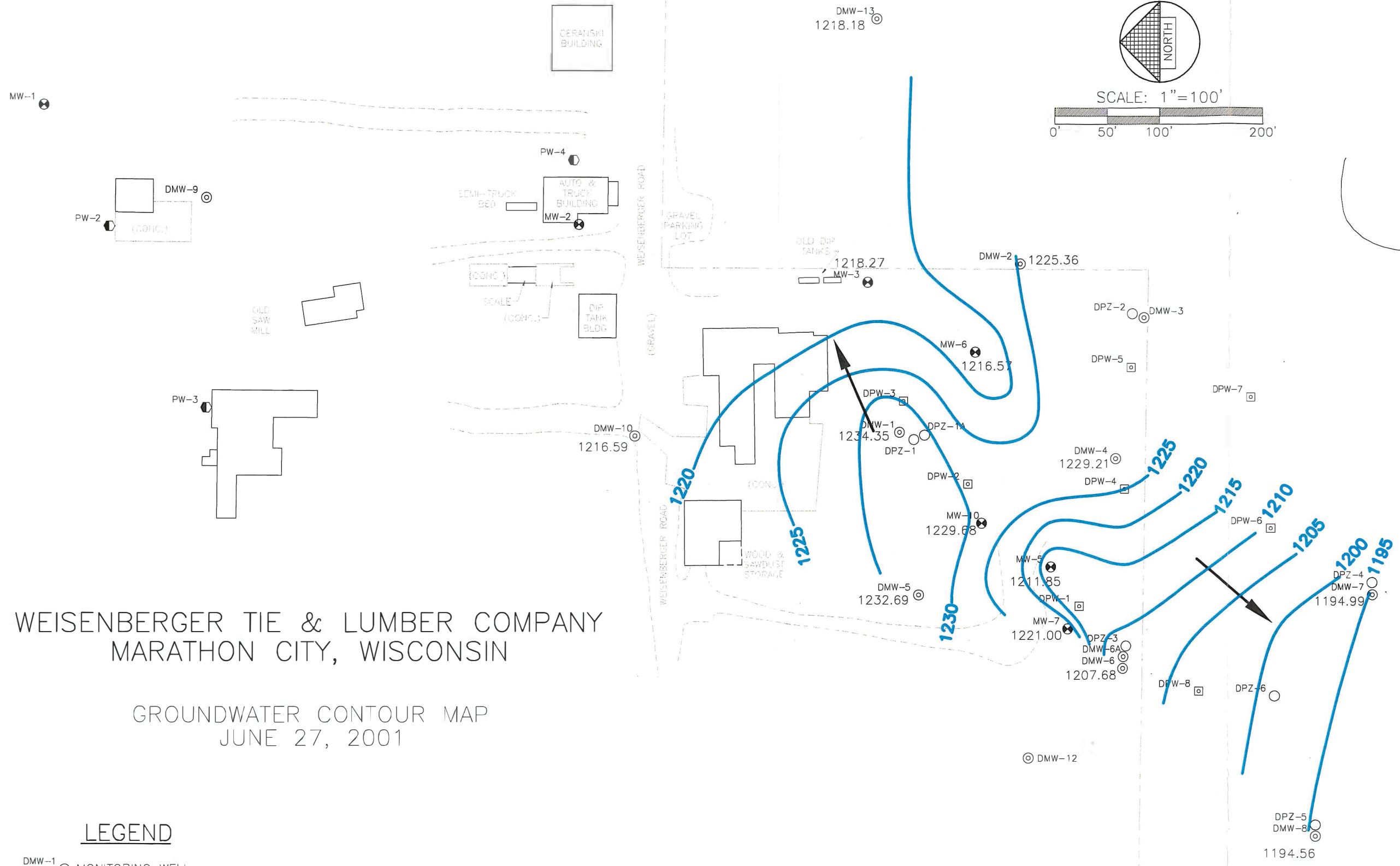
DATE SAMPLED: 6/27/01

By: CSW & BLB

METER	True / Actual	True / Actual
PH	7.00 / 7.00	4.00 / 4.00
COND	0 / 0	1413 / 1413

JOB #13551-004

WELL	PVC ELEV.	BOTTOM DEPTH	WATER DEPTH	WATER ELEV.	VOLUME BAILED GALLONS	DISSOLVED OXYGEN	PH	COND	TEMP °C	ODOR	COLOR	TURBID	COMMENTS
MW-3	1252.67	56.50	34.40	1218.27	17.0	2.5	6.70	672	16	Y	N	Y	
MW-5	1239.71	52.95	27.86	1211.85	19.0	0.6	7.11	620	15	Y	N	Y	
MW-6	1249.44	54.75	32.87	1216.57	17.0	1.9	6.33	309	14	N	N	Y	
MW-7	1237.94	30.50	16.94	1221.00	11.0	1.4	6.87	588	14	Y	N	Y	
MW-10	1242.28	22.15	12.60	1229.68	8.0	0.4	6.41	551	13	Y	N	Y	
MW-1	1247.51	18.05	13.16	1234.35	5.0	0.9	6.46	366	14	Y	N	Y	
MW-2	1246.65	27.30	21.29	1225.36	2.0 (Dry)	4.9	6.53	260	15	N	N	Y	
DMW-3	1241.46	27.95	DRY	NO SAMPLE									
MW-4	1241.16	19.00	11.95	1229.21	1.5 (Dry)	1.9	6.52	574	15	N	N	Y	
DMW-5	1244.86	19.00	12.17	1232.69	5.0	4.1	6.91	675	15	N	N	Y	
MW-6A	1236.89	32.84	29.21	1207.68	0.2 (Dry)	4.9	6.95	664	14	N	N	Y	
MW-7	1212.19	37.99	17.20	1194.99	6.0 (Dry)	4.2	6.57	322	18	N	N	Y	
DMW-8	1210.03	24.80	15.47	1194.56	2.0 (Dry)	3.8	6.49	279	14	N	N	Y	
MW-10	1236.68	30.44	20.09	1216.59	6.0 (Dry)	1.3	6.25	250	16	N	N	Y	
DMW-13	1232.93	54.98	14.75	1218.18	30.0	4.9	7.01	462	20	N	N	Y	
Z-1	1247.80	52.20	21.34	1226.46	22.0	0.4	7.08	874	15	Y	N	Y	
Z-1a	1248.12	110.15	24.74	1223.38	22.0 (Dry)	0.6	7.15	776	13	N	N	Y	
DPZ-2	1240.84	52.20	28.41	1212.43	18.0	2.1	6.93	371	14	N	N	Y	
Z-3	1236.65	49.18	29.78	1206.87	23.0	1.1	6.51	561	14	N	N	Y	
DPZ-4	1213.19	72.88	69.55	1143.64	0.5 (Dry)	2.4	7.11	697	17	N	N	Y	
Z-5	1209.38	67.86	15.30	1194.08	20.0 (Dry)	2.4	6.59	653	16	N	N	Y	
Z-6	1211.56	47.66	46.38	1165.18	0.2 (Dry)	1.9	6.47	207	17	N	N	Y	
DUP 1 (MW-3)						2.5	6.70	672	16	Y	N	Y	
EQUIP B1							6.09	18.3	24	N	N	N	
P B										N	N	N	
BK859 (PRIVATE WELL)							6.52	190	20	N	N	N	



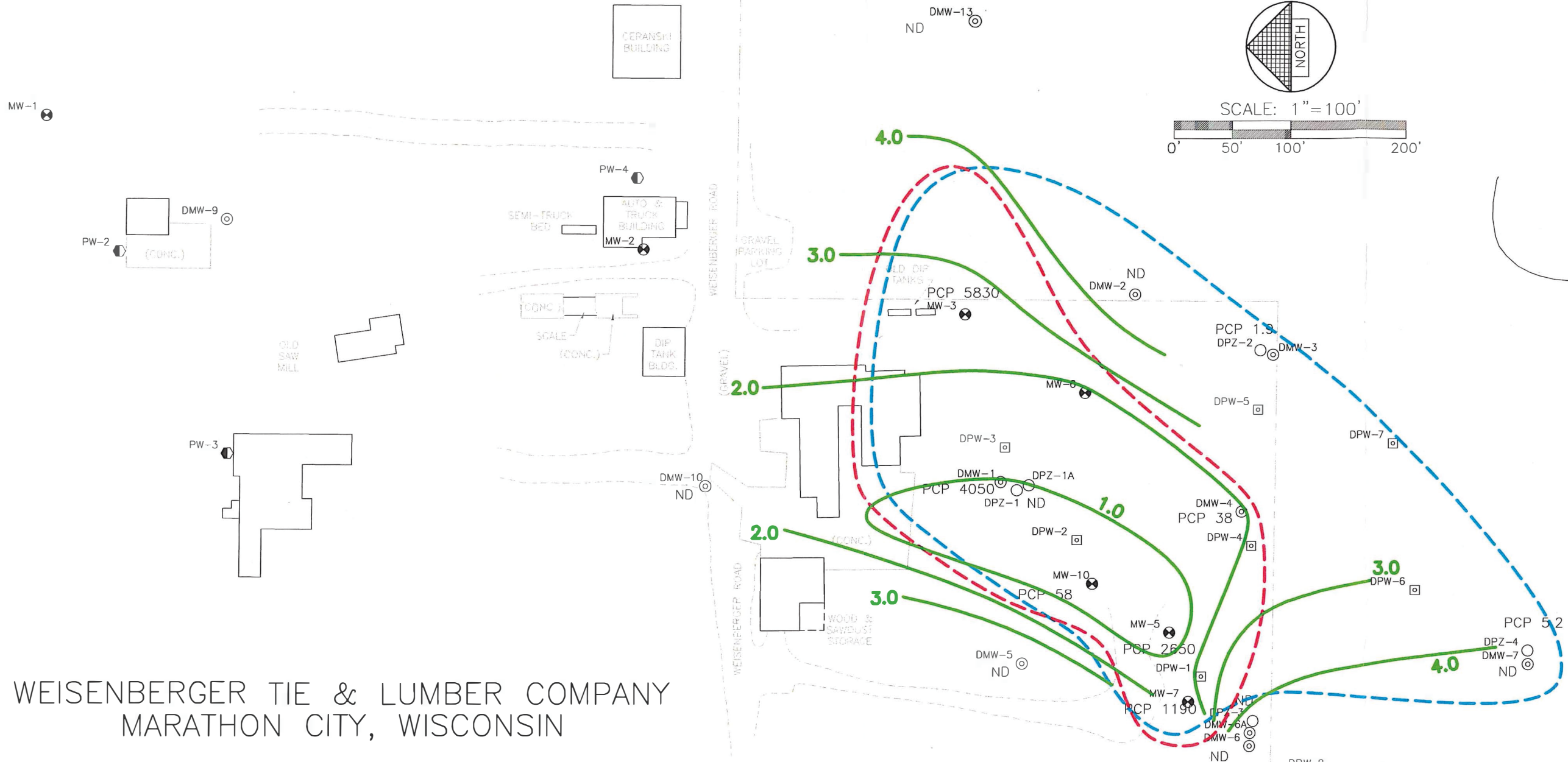
WEISENBERGER TIE & LUMBER COMPANY
MARATHON CITY, WISCONSIN

GROUNDWATER CONTOUR MAP
JUNE 27, 2001

LEGEND

- DMW-1 ⊙ MONITORING WELL
- DPZ-2 ○ PIEZOMETER
- DPW-3 □ DELTA PUMPING WELL
- MW-4 ⊙ MONITORING WELL
- PW-5 ● PRIVATE WATER SUPPLY WELL

FIGURE 1



WEISENBERGER TIE & LUMBER COMPANY
MARATHON CITY, WISCONSIN

CONTAMINANT AND OXYGEN DISTRIBUTION MAP
JUNE 27, 2001

LEGEND

- DMW-1 ⊙ MONITORING WELL
- DPZ-2 ○ PIEZOMETER
- DPW-3 □ DELTA PUMPING WELL
- MW-4 ⊕ MONITORING WELL
- PW-5 ● PRIVATE WATER SUPPLY WELL

- OXYGEN DISTRIBUTION (ppm)
- - - APPROXIMATE LIMIT OF PCP CONTAMINATION
BASED ON MONITORING WELL ANALYSIS (ppb)
- - - APPROXIMATE LIMIT OF PCP CONTAMINATION
BASED ON PIEZOMETER ANALYSIS (ppb)

FIGURE 1

LABORATORY REPORT
ROBERT E. LEE & ASSOCIATES, INC.

• MONITORING WELLS



Robert E. Lee & Associates, Inc.

Engineering, Surveying, Laboratory Services

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Wisconsin Certification Number: 405043870

JIM CAINE
ROBERT E LEE & ASSOCIATES, INC
2825 S WEBSTER AVE
GREEN BAY WI 54301-2878

Phone: (920)336-6338
Fax: (920)336-9141
Client ID: L14
Contact ID: 1859

Sample Information	Number of pages attached
Report Date: 8/23/2001	Coversheet: 1
Chain Number: 85968	Analyst generated narratives: 2
Project No: 13551002	Certificate of Analysis: 46
Project Name: WEISENBERGER TIE & LUMBER	Flag description: 1
Receive Date: 6/27/2001	Invoice: 3
Sample Date: 6/27/2001	Chain of Custody: 4
	DNR Form: 0
	Sample non-compliance Report: 0
	Subcontracted Lab Report: 40
	Miscellaneous: 0
	Total pages: 97

Attest:

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Solid sample results are reported on a dry weight basis.

ROBERT E. LEE & ASSOCIATES, INC.

CLIENT: ROBERT E LEE & ASSOCIATES, INC
PROJECT: 13551002/WEISENBERGER TIE & LUMBER
CHAIN NUMBER: 85968

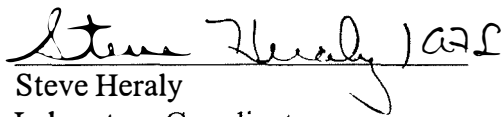
NARRATIVE

This narrative is relevant to samples MW-3, DMW-4, DMW-5, DMW-6A, DMW-10, DPZ-1, DUP-1, EQUIP B1, and TRIP B.

The samples were analyzed for petroleum volatile organic compounds following SW-846 Method 8021 and the Wisconsin Modified GRO Method.

The following is a summary of the quality control results:

1. The reported compounds were not detected in the water method blank.
2. The precision between the matrix spike recovery and matrix spike duplicate recovery was within laboratory limits for each of the reported compounds.
3. The precision between the recoveries of the water duplicate control spikes was within method limits for each of the reported compounds.
4. The matrix spike and matrix spike duplicate recoveries were within laboratory limits for each of the reported compounds except xylenes, which were below laboratory limits. These failures are believed to be caused by a matrix effect since the recoveries were low in both spiked samples. The results were accepted because the recovery for each water laboratory control spike was within method limits for all compounds.
5. The recovery for each water laboratory control spike was within method limits for each of the reported compounds.
6. The surrogate recovery for all samples was within laboratory limits.
7. The initial and final calibration check standards verified the calibration curve for each of the reported compounds.


Steve Heraly

Laboratory Coordinator

cw

ROBERT E. LEE & ASSOCIATES, INC

CLIENT: ROBERT E. LEE & ASSOCIATES, INC.
PROJECT: 13551002/WEISENBERGER TIE & LUMBER
CHAIN NUMBER: 85968

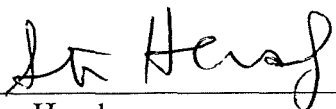
NARRATIVE

This narrative is relevant to samples DMW-10, DMW-13, DPZ-1, DPZ-1A, DPZ-2, DPZ-3, DPZ-4, DPZ-5, MW-3, MW-5, MW-6, MW-7, MW-10, DMW-1, DMW-2, DMW-4, DMW-5, DMW-6A, DMW-7 and DPZ-6.

The samples were analyzed for semi-volatile organic compounds following SW-846 Method 8270C.

Sample DMW-10 was used for the matrix spikes. The following is a summary of the quality control results:

1. The reported compounds were not detected in the method blank.
2. The precision between the matrix spike recovery and method spike recovery was within laboratory limits for each of the sixty-four compounds.
3. The matrix spike and matrix spike duplicate recoveries were within laboratory limits for each of the sixty-four compounds spiked except for naphthalene which were above laboratory limits. The failure looks to have been caused by a matrix effect since the recovery for naphthalene was above laboratory limits in both matrix spikes. The data was accepted because the method spike recovery for naphthalene was within laboratory limits.
4. The surrogate recovery was within laboratory limits for each of the six surrogates spiked in all samples.
5. The initial and final check standards verified the calibration curve for each of the reported compounds.
6. The results for pentachlorophenol in samples DPZ-1A, MW-3, MW-5 and DMW-1 were above calibration curve limits. The samples were unable to be reanalyzed within hold time due to instrument problem. The sample were analyzed over the hold time with results for pentachlorophenol; DPZ-1 was 6410 ug/L, MW-3 was 23500 ug/L, MW-5 was 4800 ug/L and DMW-1 was 9590 ug/L.



Steve Heraly
Laboratory Coordinator

JF

Robert E Lee & Associates, Inc
 2825 S Webster Ave

Green Bay, WI 54301-2878
 Project Number: 13551002

Project Name: WEISENBERGER TIE & LUMBER

Attn: Jim Caine
 Phone: (920)336-6338
 Fax: (920)336-9141
 Client ID: L14
 Chain: 85968
 Report Date: 8/23/2001

Method	Parameter Name	Result	Units	Flag	MDL	PQL	Anal Date	Analyst
Lab No	Collect Date	Sample ID						
01REL010543	6/27/2001	MW-3						
EPA 1613	Dioxin Analysis	See Attached			0	0	7/22/2001	PJK
SW-846-8021B	1,2,4-Trimethylbenzene	75	ug/L		0.23	0.7666	7/10/2001	CRW
SW-846-8021B	1,3,5-Trimethylbenzene	15	ug/L		0.21	0.7	7/10/2001	CRW
SW-846-8021B	Benzene	3.6	ug/L		0.21	0.7	7/10/2001	CRW
SW-846-8021B	Ethylbenzene	11	ug/L		0.23	0.7666	7/10/2001	CRW
SW-846-8021B	Fluorobenzene-Surrogate	98	% Rec		0	0	7/10/2001	CRW
SW-846-8021B	Methyl-tertiary-butyl ether	<0.091	ug/L		0.091	0.3033	7/10/2001	CRW
SW-846-8021B	Toluene	2.0	ug/L		0.22	0.7333	7/10/2001	CRW
SW-846-8021B	Xylenes-Total	69	ug/L		0.44	1.4666	7/10/2001	CRW
SW-846-8270C	Extraction Date	Complete			0	0	7/03/2001	JF
SW-846-8270C	1,2,4,5-Tetrachlorobenzene	<3.0	ug/L		3	10	8/09/2001	JF
SW-846-8270C	1,2,4-Trichlorobenzene	<3.6	ug/L		3.6	12	8/09/2001	JF
SW-846-8270C	1,2-Dichlorobenzene	<3.2	ug/L		3.2	10.666	8/09/2001	JF
SW-846-8270C	1,2-Diphenylhydrazine	<1.3	ug/L		1.34	4.4666	8/09/2001	JF
SW-846-8270C	1,3-Dichlorobenzene	<3.6	ug/L		3.6	12	8/09/2001	JF
SW-846-8270C	1,4-Dichlorobenzene	<3.4	ug/L		3.4	11.333	8/09/2001	JF
SW-846-8270C	1-Chloronaphthalene	<1.8	ug/L		1.84	6.1333	8/09/2001	JF
SW-846-8270C	1-Naphthylamine	<2.2	ug/L		2.2	7.3333	8/09/2001	JF
SW-846-8270C	2,3,4,6-Tetrachlorophenol	552	ug/L		1.36	4.5333	8/09/2001	JF
SW-846-8270C	2,4,5-Trichlorophenol	<1.6	ug/L		1.64	5.4666	8/09/2001	JF
SW-846-8270C	2,4,6-Tribromophenol - Surrogate	83	% Rec		0	0	8/09/2001	JF
SW-846-8270C	2,4,6-Trichlorophenol	<1.6	ug/L		1.62	5.4	8/09/2001	JF
SW-846-8270C	2,4-Dichlorophenol	<1.5	ug/L		1.54	5.1333	8/09/2001	JF
SW-846-8270C	2,4-Dimethylphenol	<2.8	ug/L		2.8	9.3333	8/09/2001	JF
SW-846-8270C	2,4-Dinitrophenol	<1.8	ug/L		1.78	5.9333	8/09/2001	JF
SW-846-8270C	2,4-Dinitrotoluene	<1.6	ug/L		1.58	5.2666	8/09/2001	JF
SW-846-8270C	2,6-Dichlorophenol	<1.5	ug/L		1.52	5.0666	8/09/2001	JF
SW-846-8270C	2,6-Dinitrotoluene	<1.6	ug/L		1.6	5.3333	8/09/2001	JF
SW-846-8270C	2-Chloronaphthalene	<1.6	ug/L		1.56	5.2	8/09/2001	JF
SW-846-8270C	2-Chlorophenol	<1.2	ug/L		1.2	4	8/09/2001	JF
SW-846-8270C	2-Fluorobiphenyl - Surrogate	69	% Rec		0	0	8/09/2001	JF
SW-846-8270C	2-Fluorophenol - Surrogate	59	% Rec		0	0	8/09/2001	JF
SW-846-8270C	2-Methylnaphthalene	203	ug/L		1.6	5.3333	8/09/2001	JF
SW-846-8270C	2-Methylphenol	<1.7	ug/L		1.7	5.6666	8/09/2001	JF
SW-846-8270C	2-Naphthylamine	<2.2	ug/L		2.2	7.3333	8/09/2001	JF
SW-846-8270C	2-Nitroaniline	<1.6	ug/L		1.64	5.4666	8/09/2001	JF
SW-846-8270C	2-Nitrophenol	<1.3	ug/L		1.28	4.2666	8/09/2001	JF
SW-846-8270C	2-Picoline	<3.0	ug/L		3	10	8/09/2001	JF
SW-846-8270C	3 & 4-Methylphenol	<1.8	ug/L		1.76	5.8666	8/09/2001	JF
SW-846-8270C	3,3'-Dichlorobenzidine	<0.84	ug/L		0.84	2.8	8/09/2001	JF
SW-846-8270C	3-Methylcholanthrene	<1.9	ug/L		1.94	6.4666	8/09/2001	JF
SW-846-8270C	3-Nitroaniline	<2.2	ug/L		2.2	7.3333	8/09/2001	JF
SW-846-8270C	4,6-Dinitro-2-methylphenol	<2.6	ug/L		2.6	8.6666	8/09/2001	JF
SW-846-8270C	4-Aminobiphenyl	<1.5	ug/L		1.5	5	8/09/2001	JF
SW-846-8270C	4-Bromophenyl phenyl ether	<1.2	ug/L		1.2	4	8/09/2001	JF

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Client ID: L14

Chain: 85968

Project Name: WEISENBERGER TIE & LUMBER

Report Date: 8/23/2001

Method	Parameter Name	Result	Units	Flag	MDL	PQL	Anl Date	Analyst
Lab No	Collect Date	Sample ID						
SW-846-8270C	4-Chloro-3-methylphenol	<1.7	ug/L		1.7	5.6666	8/09/2001	JF
SW-846-8270C	4-Chloroaniline	<1.3	ug/L		1.32	4.4	8/09/2001	JF
SW-846-8270C	4-Chlorophenyl phenyl ether	<1.2	ug/L		1.24	4.1333	8/09/2001	JF
SW-846-8270C	4-Nitroaniline	<1.3	ug/L		1.32	4.4	8/09/2001	JF
SW-846-8270C	4-Nitrophenol	<1.9	ug/L		1.86	6.2	8/09/2001	JF
SW-846-8270C	7,12-Dimethylbenz(a)anthracene	<2.0	ug/L		2	6.6666	8/09/2001	JF
SW-846-8270C	Acenaphthene	7.3	ug/L		1.12	3.7333	8/09/2001	JF
SW-846-8270C	Acenaphthylene	<1.5	ug/L		1.46	4.8666	8/09/2001	JF
SW-846-8270C	Acetophenone	<1.6	ug/L		1.6	5.3333	8/09/2001	JF
SW-846-8270C	Aniline	<1.7	ug/L		1.74	5.8	8/09/2001	JF
SW-846-8270C	Anthracene	9.5	ug/L		1.3	4.3333	8/09/2001	JF
SW-846-8270C	Benzidine	<2.8	ug/L		2.8	9.3333	8/09/2001	JF
SW-846-8270C	Benzo(a)anthracene	<1.4	ug/L		1.42	4.7333	8/09/2001	JF
SW-846-8270C	Benzo(a)pyrene	<1.1	ug/L		1.1	3.6666	8/09/2001	JF
SW-846-8270C	Benzo(b)fluoranthene	<1.3	ug/L		1.28	4.2666	8/09/2001	JF
SW-846-8270C	Benzo(g,h,i)perylene	<1.3	ug/L		1.32	4.4	8/09/2001	JF
SW-846-8270C	Benzo(k)fluoranthene	<1.6	ug/L		1.58	5.2666	8/09/2001	JF
SW-846-8270C	Benzyl alcohol	<1.8	ug/L		1.8	6	8/09/2001	JF
SW-846-8270C	bis(2-Chloroethoxy)methane	<1.5	ug/L		1.48	4.9333	8/09/2001	JF
SW-846-8270C	bis(2-Chloroethyl)ether	<1.9	ug/L		1.9	6.3333	8/09/2001	JF
SW-846-8270C	bis(2-Chloroisopropyl)ether	<1.5	ug/L		1.54	5.1333	8/09/2001	JF
SW-846-8270C	bis(2-Ethylhexyl)phthalate	11	ug/L		4.6	15.333	8/09/2001	JF
SW-846-8270C	Butylbenzylphthalate	<1.8	ug/L		1.8	6	8/09/2001	JF
SW-846-8270C	Chrysene	<1.6	ug/L		1.56	5.2	8/09/2001	JF
SW-846-8270C	Di-n-butylphthalate	<1.5	ug/L		1.52	5.0666	8/09/2001	JF
SW-846-8270C	Di-n-octylphthalate	<1.5	ug/L		1.52	5.0666	8/09/2001	JF
SW-846-8270C	Dibenz(a,j)acridine	<1.4	ug/L		1.44	4.8	8/09/2001	JF
SW-846-8270C	Dibenzo(a,h)anthracene	<1.5	ug/L		1.5	5	8/09/2001	JF
SW-846-8270C	Dibenzofuran	6.8	ug/L		1.28	4.2666	8/09/2001	JF
SW-846-8270C	Diethylphthalate	<1.3	ug/L		1.26	4.2	8/09/2001	JF
SW-846-8270C	Dimethylphthalate	<1.5	ug/L		1.52	5.0666	8/09/2001	JF
SW-846-8270C	Diphenylamine	<2.6	ug/L		2.6	8.6666	8/09/2001	JF
SW-846-8270C	Ethyl methanesulfonate	<1.7	ug/L		1.7	5.6666	8/09/2001	JF
SW-846-8270C	Fluoranthene	<2.0	ug/L		1.98	6.6	8/09/2001	JF
SW-846-8270C	Fluorene	11	ug/L		1.2	4	8/09/2001	JF
SW-846-8270C	Hexachlorobenzene	<1.6	ug/L		1.58	5.2666	8/09/2001	JF
SW-846-8270C	Hexachlorobutadiene	<7.6	ug/L		7.6	25.333	8/09/2001	JF
SW-846-8270C	Hexachlorocyclopentadiene	<5.4	ug/L		5.4	18	8/09/2001	JF
SW-846-8270C	Hexachloroethane	<5.2	ug/L		5.2	17.333	8/09/2001	JF
SW-846-8270C	Indeno(1,2,3-cd)pyrene	<1.4	ug/L		1.44	4.8	8/09/2001	JF
SW-846-8270C	Isophorone	37	ug/L		1.42	4.7333	8/09/2001	JF
SW-846-8270C	Methyl methanesulfonate	<1.1	ug/L		1.06	3.5333	8/09/2001	JF
SW-846-8270C	n-Nitrosodi-n-butylamine	<1.5	ug/L		1.46	4.8666	8/09/2001	JF
SW-846-8270C	n-Nitrosodi-n-propylamine	<1.4	ug/L		1.4	4.6666	8/09/2001	JF
SW-846-8270C	n-Nitrosodimethylamine	<1.9	ug/L		1.88	6.2666	8/09/2001	JF
SW-846-8270C	n-Nitrosodiphenylamine	<2.6	ug/L		2.6	8.6666	8/09/2001	JF

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Project Name: WEISENBERGER TIE & LUMBER

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Method	Parameter Name	Result	Units	Flag	MDL	PQL	Anl Date	Analyst
Lab No.	Collect Date	Sample ID						
SW-846-8270C	n-Nitrosopiperidine	<1.4	ug/L		1.4	4.6666	8/09/2001	JF
SW-846-8270C	Naphthalene	60	ug/L		1.38	4.6	8/09/2001	JF
SW-846-8270C	Nitrobenzene	<1.4	ug/L		1.42	4.7333	8/09/2001	JF
SW-846-8270C	Nitrobenzene-d5 - Surrogate	69	% Rec		0	0	8/09/2001	JF
SW-846-8270C	p-Dimethylaminoazobenzene	<2.0	ug/L		2	6.6666	8/09/2001	JF
SW-846-8270C	Pentachlorobenzene	<2.2	ug/L		2.2	7.3333	8/09/2001	JF
SW-846-8270C	Pentachloronitrobenzene	<1.5	ug/L		1.54	5.1333	8/09/2001	JF
SW-846-8270C	Pentachlorophenol	5830	ug/L		1.72	5.7333	8/09/2001	JF
SW-846-8270C	Phenacetin	<1.8	ug/L		1.78	5.9333	8/09/2001	JF
SW-846-8270C	Phenanthrene	<1.4	ug/L		1.42	4.7333	8/09/2001	JF
SW-846-8270C	Phenol	<1.1	ug/L		1.1	3.6666	8/09/2001	JF
SW-846-8270C	Phenol-d5 - Surrogate	51	% Rec		0	0	8/09/2001	JF
SW-846-8270C	Pronamide	<1.7	ug/L		1.68	5.6	8/09/2001	JF
SW-846-8270C	Pyrene	<2.2	ug/L		2.2	7.3333	8/09/2001	JF
SW-846-8270C	Pyridine	<3.0	ug/L		3	10	8/09/2001	JF
SW-846-8270C	Terphenyl-d14 - Surrogate	59	% Rec		0	0	8/09/2001	JF
01REL010544	6/27/2001	MW-5						
EPA 1613	Dioxin Analysis	See Attached			0	0	7/22/2001	PJK
SW-846-8270C	Extraction Date	Complete			0	0	7/03/2001	JF
SW-846-8270C	1,2,4,5-Tetrachlorobenzene	<3.0	ug/L		3	10	8/09/2001	JF
SW-846-8270C	1,2,4-Trichlorobenzene	<3.6	ug/L		3.6	12	8/09/2001	JF
SW-846-8270C	1,2-Dichlorobenzene	<3.2	ug/L		3.2	10.666	8/09/2001	JF
SW-846-8270C	1,2-Diphenylhydrazine	<1.3	ug/L		1.34	4.4666	8/09/2001	JF
SW-846-8270C	1,3-Dichlorobenzene	<3.6	ug/L		3.6	12	8/09/2001	JF
SW-846-8270C	1,4-Dichlorobenzene	<3.4	ug/L		3.4	11.333	8/09/2001	JF
SW-846-8270C	1-Chloronaphthalene	<1.8	ug/L		1.84	6.1333	8/09/2001	JF
SW-846-8270C	1-Naphthylamine	<2.2	ug/L		2.2	7.3333	8/09/2001	JF
SW-846-8270C	2,3,4,6-Tetrachlorophenol	56	ug/L		1.36	4.5333	8/09/2001	JF
SW-846-8270C	2,4,5-Trichlorophenol	<1.6	ug/L		1.64	5.4666	8/09/2001	JF
SW-846-8270C	2,4,6-Tribromophenol - Surrogate	78	% Rec		0	0	8/09/2001	JF
SW-846-8270C	2,4,6-Trichlorophenol	<1.6	ug/L		1.62	5.4	8/09/2001	JF
SW-846-8270C	2,4-Dichlorophenol	<1.5	ug/L		1.54	5.1333	8/09/2001	JF
SW-846-8270C	2,4-Dimethylphenol	<2.8	ug/L		2.8	9.3333	8/09/2001	JF
SW-846-8270C	2,4-Dinitrophenol	<1.8	ug/L		1.78	5.9333	8/09/2001	JF
SW-846-8270C	2,4-Dinitrotoluene	<1.6	ug/L		1.58	5.2666	8/09/2001	JF
SW-846-8270C	2,6-Dichlorophenol	<1.5	ug/L		1.52	5.0666	8/09/2001	JF
SW-846-8270C	2,6-Dinitrotoluene	<1.6	ug/L		1.6	5.3333	8/09/2001	JF
SW-846-8270C	2-Chloronaphthalene	<1.6	ug/L		1.56	5.2	8/09/2001	JF
SW-846-8270C	2-Chlorophenol	<1.2	ug/L		1.2	4	8/09/2001	JF
SW-846-8270C	2-Fluorobiphenyl - Surrogate	67	% Rec		0	0	8/09/2001	JF
SW-846-8270C	2-Fluorophenol - Surrogate	53	% Rec		0	0	8/09/2001	JF
SW-846-8270C	2-Methylnaphthalene	89	ug/L		1.6	5.3333	8/09/2001	JF
SW-846-8270C	2-Methylphenol	<1.7	ug/L		1.7	5.6666	8/09/2001	JF
SW-846-8270C	2-Naphthylamine	<2.2	ug/L		2.2	7.3333	8/09/2001	JF
SW-846-8270C	2-Nitroaniline	<1.6	ug/L		1.64	5.4666	8/09/2001	JF
SW-846-8270C	2-Nitrophenol	<1.3	ug/L		1.28	4.2666	8/09/2001	JF

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Method	Parameter Name	Result	Units	Flag	MDL	PQL	Anl Date	Analyst
Lab No	Collect Date	Sample ID						
SW-846-8270C	2-Picoline	<3.0	ug/L		3	10	8/09/2001	JF
SW-846-8270C	3 & 4-Methylphenol	<1.8	ug/L		1.76	5.8666	8/09/2001	JF
SW-846-8270C	3,3'-Dichlorobenzidine	<0.84	ug/L		0.84	2.8	8/09/2001	JF
SW-846-8270C	3-Methylcholanthrene	<1.9	ug/L		1.94	6.4666	8/09/2001	JF
SW-846-8270C	3-Nitroaniline	<2.2	ug/L		2.2	7.3333	8/09/2001	JF
SW-846-8270C	4,6-Dinitro-2-methylphenol	<2.6	ug/L		2.6	8.6666	8/09/2001	JF
SW-846-8270C	4-Aminobiphenyl	<1.5	ug/L		1.5	5	8/09/2001	JF
SW-846-8270C	4-Bromophenyl phenyl ether	<1.2	ug/L		1.2	4	8/09/2001	JF
SW-846-8270C	4-Chloro-3-methylphenol	<1.7	ug/L		1.7	5.6666	8/09/2001	JF
SW-846-8270C	4-Chloroaniline	<1.3	ug/L		1.32	4.4	8/09/2001	JF
SW-846-8270C	4-Chlorophenyl phenyl ether	<1.2	ug/L		1.24	4.1333	8/09/2001	JF
SW-846-8270C	4-Nitroaniline	<1.3	ug/L		1.32	4.4	8/09/2001	JF
SW-846-8270C	4-Nitrophenol	<1.9	ug/L		1.86	6.2	8/09/2001	JF
SW-846-8270C	7,12-Dimethylbenz(a)anthracene	<2.0	ug/L		2	6.6666	8/09/2001	JF
SW-846-8270C	Acenaphthene	<1.1	ug/L		1.12	3.7333	8/09/2001	JF
SW-846-8270C	Acenaphthylene	<1.5	ug/L		1.46	4.8666	8/09/2001	JF
SW-846-8270C	Acetophenone	<1.6	ug/L		1.6	5.3333	8/09/2001	JF
SW-846-8270C	Aniline	<1.7	ug/L		1.74	5.8	8/09/2001	JF
SW-846-8270C	Anthracene	<1.3	ug/L		1.3	4.3333	8/09/2001	JF
SW-846-8270C	Benzidine	<2.8	ug/L		2.8	9.3333	8/09/2001	JF
SW-846-8270C	Benzo(a)anthracene	<1.4	ug/L		1.42	4.7333	8/09/2001	JF
SW-846-8270C	Benzo(a)pyrene	<1.1	ug/L		1.1	3.6666	8/09/2001	JF
SW-846-8270C	Benzo(b)fluoranthene	<1.3	ug/L		1.28	4.2666	8/09/2001	JF
SW-846-8270C	Benzo(g,h,i)perylene	<1.3	ug/L		1.32	4.4	8/09/2001	JF
SW-846-8270C	Benzo(k)fluoranthene	<1.6	ug/L		1.58	5.2666	8/09/2001	JF
SW-846-8270C	Benzyl alcohol	<1.8	ug/L		1.8	6	8/09/2001	JF
SW-846-8270C	bis(2-Chloroethoxy)methane	<1.5	ug/L		1.48	4.9333	8/09/2001	JF
SW-846-8270C	bis(2-Chloroethyl)ether	<1.9	ug/L		1.9	6.3333	8/09/2001	JF
SW-846-8270C	bis(2-Chloroisopropyl)ether	<1.5	ug/L		1.54	5.1333	8/09/2001	JF
SW-846-8270C	bis(2-Ethylhexyl)phthalate	<4.6	ug/L		4.6	15.333	8/09/2001	JF
SW-846-8270C	Butylbenzylphthalate	<1.8	ug/L		1.8	6	8/09/2001	JF
SW-846-8270C	Chrysene	<1.6	ug/L		1.56	5.2	8/09/2001	JF
SW-846-8270C	Di-n-butylphthalate	<1.5	ug/L		1.52	5.0666	8/09/2001	JF
SW-846-8270C	Di-n-octylphthalate	<1.5	ug/L		1.52	5.0666	8/09/2001	JF
SW-846-8270C	Dibenz(a,j)acridine	<1.4	ug/L		1.44	4.8	8/09/2001	JF
SW-846-8270C	Dibenzo(a,h)anthracene	<1.5	ug/L		1.5	5	8/09/2001	JF
SW-846-8270C	Dibenzofuran	4.4	ug/L		1.28	4.2666	8/09/2001	JF
SW-846-8270C	Diethylphthalate	<1.3	ug/L		1.26	4.2	8/09/2001	JF
SW-846-8270C	Dimethylphthalate	<1.5	ug/L		1.52	5.0666	8/09/2001	JF
SW-846-8270C	Diphenylamine	<2.6	ug/L		2.6	8.6666	8/09/2001	JF
SW-846-8270C	Ethyl methanesulfonate	<1.7	ug/L		1.7	5.6666	8/09/2001	JF
SW-846-8270C	Fluoranthene	<2.0	ug/L		1.98	6.6	8/09/2001	JF
SW-846-8270C	Fluorene	<1.2	ug/L		1.2	4	8/09/2001	JF
SW-846-8270C	Hexachlorobenzene	<1.6	ug/L		1.58	5.2666	8/09/2001	JF
SW-846-8270C	Hexachlorobutadiene	<7.6	ug/L		7.6	25.333	8/09/2001	JF
SW-846-8270C	Hexachlorocyclopentadiene	<5.4	ug/L		5.4	18	8/09/2001	JF

Robert E. Lee & Associates, Inc.
 Wisconsin Certification Number: 405043870
 Certificate of Analysis Report

Robert E Lee & Associates, Inc
 2825 S Webster Ave

Attn: Jim Caine
 Phone: (920)336-6338
 Fax: (920)336-9141
 Client ID: L14
 Chain: 85968
 Report Date: 8/23/2001

Green Bay, WI 54301-2878
 Project Number: 13551002
 Project Name: WEISENBERGER TIE & LUMBER

Method	Parameter Name	Result	Units	Flag	MDL	PQL	Anl Date	Analyst
Lab No.	Collect Date	Sample ID						
SW-846-8270C	Hexachloroethane	<5.2	ug/L		5.2	17.333	8/09/2001	JF
SW-846-8270C	Indeno(1,2,3-cd)pyrene	<1.4	ug/L		1.44	4.8	8/09/2001	JF
SW-846-8270C	Isophorone	<1.4	ug/L		1.42	4.7333	8/09/2001	JF
SW-846-8270C	Methyl methanesulfonate	<1.1	ug/L		1.06	3.5333	8/09/2001	JF
SW-846-8270C	n-Nitrosodi-n-butylamine	<1.5	ug/L		1.46	4.8666	8/09/2001	JF
SW-846-8270C	n-Nitrosodi-n-propylamine	<1.4	ug/L		1.4	4.6666	8/09/2001	JF
SW-846-8270C	n-Nitrosodimethylamine	<1.9	ug/L		1.88	6.2666	8/09/2001	JF
SW-846-8270C	n-Nitrosodiphenylamine	<2.6	ug/L		2.6	8.6666	8/09/2001	JF
SW-846-8270C	n-Nitrosopiperidine	<1.4	ug/L		1.4	4.6666	8/09/2001	JF
SW-846-8270C	Naphthalene	31	ug/L		1.38	4.6	8/09/2001	JF
SW-846-8270C	Nitrobenzene	<1.4	ug/L		1.42	4.7333	8/09/2001	JF
SW-846-8270C	Nitrobenzene-d5 - Surrogate	64	% Rec		0	0	8/09/2001	JF
SW-846-8270C	p-Dimethylaminoazobenzene	<2.0	ug/L		2	6.6666	8/09/2001	JF
SW-846-8270C	Pentachlorobenzene	<2.2	ug/L		2.2	7.3333	8/09/2001	JF
SW-846-8270C	Pentachloronitrobenzene	<1.5	ug/L		1.54	5.1333	8/09/2001	JF
SW-846-8270C	Pentachlorophenol	2650	ug/L		1.72	5.7333	8/09/2001	JF
SW-846-8270C	Phenacetin	<1.8	ug/L		1.78	5.9333	8/09/2001	JF
SW-846-8270C	Phenanthrene	2.5	ug/L		1.42	4.7333	8/09/2001	JF
SW-846-8270C	Phenol	<1.1	ug/L		1.1	3.6666	8/09/2001	JF
SW-846-8270C	Phenol-d5 - Surrogate	49	% Rec		0	0	8/09/2001	JF
SW-846-8270C	Pronamide	<1.7	ug/L		1.68	5.6	8/09/2001	JF
SW-846-8270C	Pyrene	<2.2	ug/L		2.2	7.3333	8/09/2001	JF
SW-846-8270C	Pyridine	<3.0	ug/L		3	10	8/09/2001	JF
SW-846-8270C	Terphenyl-d14 - Surrogate	55	% Rec		0	0	8/09/2001	JF

01REL010545 6/27/2001 MW-6

EPA 1613	Dioxin Analysis	See Attached			0	0	7/22/2001	PJK
SW-846-8270C	Extraction Date	Complete			0	0	7/03/2001	JF
SW-846-8270C	1,2,4,5-Tetrachlorobenzene	<3.0	ug/L		3	10	8/09/2001	JF
SW-846-8270C	1,2,4-Trichlorobenzene	<3.6	ug/L		3.6	12	8/09/2001	JF
SW-846-8270C	1,2-Dichlorobenzene	<3.2	ug/L		3.2	10.666	8/09/2001	JF
SW-846-8270C	1,2-Diphenylhydrazine	<1.3	ug/L		1.34	4.4666	8/09/2001	JF
SW-846-8270C	1,3-Dichlorobenzene	<3.6	ug/L		3.6	12	8/09/2001	JF
SW-846-8270C	1,4-Dichlorobenzene	<3.4	ug/L		3.4	11.333	8/09/2001	JF
SW-846-8270C	1-Chloronaphthalene	<1.8	ug/L		1.84	6.1333	8/09/2001	JF
SW-846-8270C	1-Naphthylamine	<2.2	ug/L		2.2	7.3333	8/09/2001	JF
SW-846-8270C	2,3,4,6-Tetrachlorophenol	2.64	ug/L	13	1.36	4.5333	8/09/2001	JF
SW-846-8270C	2,4,5-Trichlorophenol	<1.6	ug/L		1.64	5.4666	8/09/2001	JF
SW-846-8270C	2,4,6-Tribromophenol - Surrogate	72	% Rec		0	0	8/09/2001	JF
SW-846-8270C	2,4,6-Trichlorophenol	<1.6	ug/L		1.62	5.4	8/09/2001	JF
SW-846-8270C	2,4-Dichlorophenol	<1.5	ug/L		1.54	5.1333	8/09/2001	JF
SW-846-8270C	2,4-Dimethylphenol	<2.8	ug/L		2.8	9.3333	8/09/2001	JF
SW-846-8270C	2,4-Dinitrophenol	<1.8	ug/L		1.78	5.9333	8/09/2001	JF
SW-846-8270C	2,4-Dinitrotoluene	<1.6	ug/L		1.58	5.2666	8/09/2001	JF
SW-846-8270C	2,6-Dichlorophenol	<1.5	ug/L		1.52	5.0666	8/09/2001	JF
SW-846-8270C	2,6-Dinitrotoluene	<1.6	ug/L		1.6	5.3333	8/09/2001	JF
SW-846-8270C	2-Chloronaphthalene	<1.6	ug/L		1.56	5.2	8/09/2001	JF

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Green Bay, WI 54301-2878

Project Number: 13551002

Client ID: L14

Chain: 85968

Project Name: WEISENBERGER TIE & LUMBER

Report Date: 8/23/2001

Method	Parameter Name	Result	Units	Flag	MDL	PQL	Anl Date	Analyst
Lab No	Collect Date	Sample ID						
SW-846-8270C	2-Chlorophenol	<1.2	ug/L		1.2	4	8/09/2001	JF
SW-846-8270C	2-Fluorobiphenyl - Surrogate	61	% Rec		0	0	8/09/2001	JF
SW-846-8270C	2-Fluorophenol - Surrogate	50	% Rec		0	0	8/09/2001	JF
SW-846-8270C	2-Methylnaphthalene	<1.6	ug/L		1.6	5.3333	8/09/2001	JF
SW-846-8270C	2-Methylphenol	<1.7	ug/L		1.7	5.6666	8/09/2001	JF
SW-846-8270C	2-Naphthylamine	<2.2	ug/L		2.2	7.3333	8/09/2001	JF
SW-846-8270C	2-Nitroaniline	<1.6	ug/L		1.64	5.4666	8/09/2001	JF
SW-846-8270C	2-Nitrophenol	<1.3	ug/L		1.28	4.2666	8/09/2001	JF
SW-846-8270C	2-Picoline	<3.0	ug/L		3	10	8/09/2001	JF
SW-846-8270C	3 & 4-Methylphenol	<1.8	ug/L		1.76	5.8666	8/09/2001	JF
SW-846-8270C	3,3'-Dichlorobenzidine	<0.84	ug/L		0.84	2.8	8/09/2001	JF
SW-846-8270C	3-Methylcholanthrene	<1.9	ug/L		1.94	6.4666	8/09/2001	JF
SW-846-8270C	3-Nitroaniline	<2.2	ug/L		2.2	7.3333	8/09/2001	JF
SW-846-8270C	4,6-Dinitro-2-methylphenol	<2.6	ug/L		2.6	8.6666	8/09/2001	JF
SW-846-8270C	4-Aminobiphenyl	<1.5	ug/L		1.5	5	8/09/2001	JF
SW-846-8270C	4-Bromophenyl phenyl ether	<1.2	ug/L		1.2	4	8/09/2001	JF
SW-846-8270C	4-Chloro-3-methylphenol	<1.7	ug/L		1.7	5.6666	8/09/2001	JF
SW-846-8270C	4-Chloroaniline	<1.3	ug/L		1.32	4.4	8/09/2001	JF
SW-846-8270C	4-Chlorophenyl phenyl ether	<1.2	ug/L		1.24	4.1333	8/09/2001	JF
SW-846-8270C	4-Nitroaniline	<1.3	ug/L		1.32	4.4	8/09/2001	JF
SW-846-8270C	4-Nitrophenol	<1.9	ug/L		1.86	6.2	8/09/2001	JF
SW-846-8270C	7,12-Dimethylbenz(a)anthracene	<2.0	ug/L		2	6.6666	8/09/2001	JF
SW-846-8270C	Acenaphthene	<1.1	ug/L		1.12	3.7333	8/09/2001	JF
SW-846-8270C	Acenaphthylene	<1.5	ug/L		1.46	4.8666	8/09/2001	JF
SW-846-8270C	Acetophenone	<1.6	ug/L		1.6	5.3333	8/09/2001	JF
SW-846-8270C	Aniline	<1.7	ug/L		1.74	5.8	8/09/2001	JF
SW-846-8270C	Anthracene	<1.3	ug/L		1.3	4.3333	8/09/2001	JF
SW-846-8270C	Benzenzidine	<2.8	ug/L		2.8	9.3333	8/09/2001	JF
SW-846-8270C	Benzo(a)anthracene	<1.4	ug/L		1.42	4.7333	8/09/2001	JF
SW-846-8270C	Benzo(a)pyrene	<1.1	ug/L		1.1	3.6666	8/09/2001	JF
SW-846-8270C	Benzo(b)fluoranthene	<1.3	ug/L		1.28	4.2666	8/09/2001	JF
SW-846-8270C	Benzo(g,h,i)perylene	<1.3	ug/L		1.32	4.4	8/09/2001	JF
SW-846-8270C	Benzo(k)fluoranthene	<1.6	ug/L		1.58	5.2666	8/09/2001	JF
SW-846-8270C	Benzyl alcohol	<1.8	ug/L		1.8	6	8/09/2001	JF
SW-846-8270C	bis(2-Chloroethoxy)methane	<1.5	ug/L		1.48	4.9333	8/09/2001	JF
SW-846-8270C	bis(2-Chloroethyl)ether	<1.9	ug/L		1.9	6.3333	8/09/2001	JF
SW-846-8270C	bis(2-Chloroisopropyl)ether	<1.5	ug/L		1.54	5.1333	8/09/2001	JF
SW-846-8270C	bis(2-Ethylhexyl)phthalate	<4.6	ug/L		4.6	15.333	8/09/2001	JF
SW-846-8270C	Butylbenzylphthalate	<1.8	ug/L		1.8	6	8/09/2001	JF
SW-846-8270C	Chrysene	<1.6	ug/L		1.56	5.2	8/09/2001	JF
SW-846-8270C	Di-n-butylphthalate	<1.5	ug/L		1.52	5.0666	8/09/2001	JF
SW-846-8270C	Di-n-octylphthalate	<1.5	ug/L		1.52	5.0666	8/09/2001	JF
SW-846-8270C	Dibenz(a,j)acridine	<1.4	ug/L		1.44	4.8	8/09/2001	JF
SW-846-8270C	Dibenzo(a,h)anthracene	<1.5	ug/L		1.5	5	8/09/2001	JF
SW-846-8270C	Dibenzofuran	<1.3	ug/L		1.28	4.2666	8/09/2001	JF
SW-846-8270C	Diethylphthalate	<1.3	ug/L		1.26	4.2	8/09/2001	JF

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2825 S Webster Ave

Green Bay, WI 54301-2878
Project Number: 13551002
Project Name: WEISENBERGER TIE & LUMBER

Attn: Jim Caine
Phone: (920)336-6338
Fax: (920)336-9141
Client ID: L14
Chain: 85968
Report Date: 8/23/2001

Method	Parameter Name	Result	Units	Flag	MDL	PQL	Anal Date	Analyst
Lab No	Collect Date	Sample ID						
SW-846-8270C	Dimethylphthalate	<1.5	ug/L		1.52	5.0666	8/09/2001	JF
SW-846-8270C	Diphenylamine	<2.6	ug/L		2.6	8.6666	8/09/2001	JF
SW-846-8270C	Ethyl methanesulfonate	<1.7	ug/L		1.7	5.6666	8/09/2001	JF
SW-846-8270C	Fluoranthene	<2.0	ug/L		1.98	6.6	8/09/2001	JF
SW-846-8270C	Fluorene	<1.2	ug/L		1.2	4	8/09/2001	JF
SW-846-8270C	Hexachlorobenzene	<1.6	ug/L		1.58	5.2666	8/09/2001	JF
SW-846-8270C	Hexachlorobutadiene	<7.6	ug/L		7.6	25.333	8/09/2001	JF
SW-846-8270C	Hexachlorocyclopentadiene	<5.4	ug/L		5.4	18	8/09/2001	JF
SW-846-8270C	Hexachloroethane	<5.2	ug/L		5.2	17.333	8/09/2001	JF
SW-846-8270C	Indeno(1,2,3-cd)pyrene	<1.4	ug/L		1.44	4.8	8/09/2001	JF
SW-846-8270C	Isophorone	<1.4	ug/L		1.42	4.7333	8/09/2001	JF
SW-846-8270C	Methyl methanesulfonate	<1.1	ug/L		1.06	3.5333	8/09/2001	JF
SW-846-8270C	n-Nitrosodi-n-butylamine	<1.5	ug/L		1.46	4.8666	8/09/2001	JF
SW-846-8270C	n-Nitrosodi-n-propylamine	<1.4	ug/L		1.4	4.6666	8/09/2001	JF
SW-846-8270C	n-Nitrosodimethylamine	<1.9	ug/L		1.88	6.2666	8/09/2001	JF
SW-846-8270C	n-Nitrosodiphenylamine	<2.6	ug/L		2.6	8.6666	8/09/2001	JF
SW-846-8270C	n-Nitrosopiperidine	<1.4	ug/L		1.4	4.6666	8/09/2001	JF
SW-846-8270C	Naphthalene	<1.4	ug/L		1.38	4.6	8/09/2001	JF
SW-846-8270C	Nitrobenzene	<1.4	ug/L		1.42	4.7333	8/09/2001	JF
SW-846-8270C	Nitrobenzene-d5 - Surrogate	58	% Rec		0	0	8/09/2001	JF
SW-846-8270C	p-Dimethylaminoazobenzene	<2.0	ug/L		2	6.6666	8/09/2001	JF
SW-846-8270C	Pentachlorobenzene	<2.2	ug/L		2.2	7.3333	8/09/2001	JF
SW-846-8270C	Pentachloronitrobenzene	<1.5	ug/L		1.54	5.1333	8/09/2001	JF
SW-846-8270C	Pentachlorophenol	85	ug/L		1.72	5.7333	8/09/2001	JF
SW-846-8270C	Phenacetin	<1.8	ug/L		1.78	5.9333	8/09/2001	JF
SW-846-8270C	Phenanthrene	<1.4	ug/L		1.42	4.7333	8/09/2001	JF
SW-846-8270C	Phenol	<1.1	ug/L		1.1	3.6666	8/09/2001	JF
SW-846-8270C	Phenol-d5 - Surrogate	43	% Rec		0	0	8/09/2001	JF
SW-846-8270C	Pronamide	<1.7	ug/L		1.68	5.6	8/09/2001	JF
SW-846-8270C	Pyrene	<2.2	ug/L		2.2	7.3333	8/09/2001	JF
SW-846-8270C	Pyridine	<3.0	ug/L		3	10	8/09/2001	JF
SW-846-8270C	Terphenyl-d14 - Surrogate	56	% Rec		0	0	8/09/2001	JF

01REL010546 6/27/2001 MW-7

EPA 1613	Dioxin Analysis	See Attached			0	0	7/22/2001	PJK
SW-846-8270C	Extraction Date	Complete			0	0	7/03/2001	JF
SW-846-8270C	1,2,4,5-Tetrachlorobenzene	<3.0	ug/L		3	10	8/09/2001	JF
SW-846-8270C	1,2,4-Trichlorobenzene	<3.6	ug/L		3.6	12	8/09/2001	JF
SW-846-8270C	1,2-Dichlorobenzene	<3.2	ug/L		3.2	10.666	8/09/2001	JF
SW-846-8270C	1,2-Diphenylhydrazine	<1.3	ug/L		1.34	4.4666	8/09/2001	JF
SW-846-8270C	1,3-Dichlorobenzene	<3.6	ug/L		3.6	12	8/09/2001	JF
SW-846-8270C	1,4-Dichlorobenzene	<3.4	ug/L		3.4	11.333	8/09/2001	JF
SW-846-8270C	1-Chloronaphthalene	<1.8	ug/L		1.84	6.1333	8/09/2001	JF
SW-846-8270C	1-Naphthylamine	<2.2	ug/L		2.2	7.3333	8/09/2001	JF
SW-846-8270C	2,3,4,6-Tetrachlorophenol	29	ug/L		1.36	4.5333	8/09/2001	JF
SW-846-8270C	2,4,5-Trichlorophenol	<1.6	ug/L		1.64	5.4666	8/09/2001	JF
SW-846-8270C	2,4,6-Tribromophenol - Surrogate	78	% Rec		0	0	8/09/2001	JF

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Chain: 85968

Report Date: 8/23/2001

Method	Parameter Name	Result	Units	Flag	MDL	PQL	Anl Date	Analyst
Lab No.	Collect Date	Sample ID						
SW-846-8270C	2,4,6-Trichlorophenol	<1.6	ug/L		1.62	5.4	8/09/2001	JF
SW-846-8270C	2,4-Dichlorophenol	<1.5	ug/L		1.54	5.1333	8/09/2001	JF
SW-846-8270C	2,4-Dimethylphenol	<2.8	ug/L		2.8	9.3333	8/09/2001	JF
SW-846-8270C	2,4-Dinitrophenol	<1.8	ug/L		1.78	5.9333	8/09/2001	JF
SW-846-8270C	2,4-Dinitrotoluene	<1.6	ug/L		1.58	5.2666	8/09/2001	JF
SW-846-8270C	2,6-Dichlorophenol	<1.5	ug/L		1.52	5.0666	8/09/2001	JF
SW-846-8270C	2,6-Dinitrotoluene	<1.6	ug/L		1.6	5.3333	8/09/2001	JF
SW-846-8270C	2-Chloronaphthalene	<1.6	ug/L		1.56	5.2	8/09/2001	JF
SW-846-8270C	2-Chlorophenol	<1.2	ug/L		1.2	4	8/09/2001	JF
SW-846-8270C	2-Fluorobiphenyl - Surrogate	64	% Rec		0	0	8/09/2001	JF
SW-846-8270C	2-Fluorophenol - Surrogate	47	% Rec		0	0	8/09/2001	JF
SW-846-8270C	2-Methylnaphthalene	275	ug/L		1.6	5.3333	8/09/2001	JF
SW-846-8270C	2-Methylphenol	<1.7	ug/L		1.7	5.6666	8/09/2001	JF
SW-846-8270C	2-Naphthylamine	<2.2	ug/L		2.2	7.3333	8/09/2001	JF
SW-846-8270C	2-Nitroaniline	<1.6	ug/L		1.64	5.4666	8/09/2001	JF
SW-846-8270C	2-Nitrophenol	<1.3	ug/L		1.28	4.2666	8/09/2001	JF
SW-846-8270C	2-Picoline	<3.0	ug/L		3	10	8/09/2001	JF
SW-846-8270C	3 & 4-Methylphenol	<1.8	ug/L		1.76	5.8666	8/09/2001	JF
SW-846-8270C	3,3'-Dichlorobenzidine	<0.84	ug/L		0.84	2.8	8/09/2001	JF
SW-846-8270C	3-Methylcholanthrene	<1.9	ug/L		1.94	6.4666	8/09/2001	JF
SW-846-8270C	3-Nitroaniline	<2.2	ug/L		2.2	7.3333	8/09/2001	JF
SW-846-8270C	4,6-Dinitro-2-methylphenol	<2.6	ug/L		2.6	8.6666	8/09/2001	JF
SW-846-8270C	4-Aminobiphenyl	<1.5	ug/L		1.5	5	8/09/2001	JF
SW-846-8270C	4-Bromophenyl phenyl ether	<1.2	ug/L		1.2	4	8/09/2001	JF
SW-846-8270C	4-Chloro-3-methylphenol	<1.7	ug/L		1.7	5.6666	8/09/2001	JF
SW-846-8270C	4-Chloroaniline	<1.3	ug/L		1.32	4.4	8/09/2001	JF
SW-846-8270C	4-Chlorophenyl phenyl ether	<1.2	ug/L		1.24	4.1333	8/09/2001	JF
SW-846-8270C	4-Nitroaniline	<1.3	ug/L		1.32	4.4	8/09/2001	JF
SW-846-8270C	4-Nitrophenol	<1.9	ug/L		1.86	6.2	8/09/2001	JF
SW-846-8270C	7,12-Dimethylbenz(a)anthracene	<2.0	ug/L		2	6.6666	8/09/2001	JF
SW-846-8270C	Acenaphthene	7.7	ug/L		1.12	3.7333	8/09/2001	JF
SW-846-8270C	Acenaphthylene	<1.5	ug/L		1.46	4.8666	8/09/2001	JF
SW-846-8270C	Acetophenone	<1.6	ug/L		1.6	5.3333	8/09/2001	JF
SW-846-8270C	Aniline	<1.7	ug/L		1.74	5.8	8/09/2001	JF
SW-846-8270C	Anthracene	<1.3	ug/L		1.3	4.3333	8/09/2001	JF
SW-846-8270C	Benzidine	<2.8	ug/L		2.8	9.3333	8/09/2001	JF
SW-846-8270C	Benzo(a)anthracene	<1.4	ug/L		1.42	4.7333	8/09/2001	JF
SW-846-8270C	Benzo(a)pyrene	<1.1	ug/L		1.1	3.6666	8/09/2001	JF
SW-846-8270C	Benzo(b)fluoranthene	<1.3	ug/L		1.28	4.2666	8/09/2001	JF
SW-846-8270C	Benzo(g,h,i)perylene	<1.3	ug/L		1.32	4.4	8/09/2001	JF
SW-846-8270C	Benzo(k)fluoranthene	<1.6	ug/L		1.58	5.2666	8/09/2001	JF
SW-846-8270C	Benzyl alcohol	<1.8	ug/L		1.8	6	8/09/2001	JF
SW-846-8270C	bis(2-Chloroethoxy)methane	<1.5	ug/L		1.48	4.9333	8/09/2001	JF
SW-846-8270C	bis(2-Chloroethyl)ether	<1.9	ug/L		1.9	6.3333	8/09/2001	JF
SW-846-8270C	bis(2-Chloroisopropyl)ether	<1.5	ug/L		1.54	5.1333	8/09/2001	JF
SW-846-8270C	bis(2-Ethylhexyl)phthalate	<4.6	ug/L		4.6	15.333	8/09/2001	JF

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Project Number: 13551002

Project Name: WEISENBERGER TIE & LUMBER

Attn: Jim Caine

Phone: (920)336-6338

Fax: (920)336-9141

Client ID: L14

Chain: 85968

Report Date: 8/23/2001

Method	Parameter Name	Result	Units	Flag	MDL	PQL	Anl Date	Analyst
Lab No	Collect Date	Sample ID						
SW-846-8270C	Butylbenzylphthalate	<1.8	ug/L		1.8	6	8/09/2001	JF
SW-846-8270C	Chrysene	<1.6	ug/L		1.56	5.2	8/09/2001	JF
SW-846-8270C	Di-n-butylphthalate	<1.5	ug/L		1.52	5.0666	8/09/2001	JF
SW-846-8270C	Di-n-octylphthalate	<1.5	ug/L		1.52	5.0666	8/09/2001	JF
SW-846-8270C	Dibenz(a,j)acridine	<1.4	ug/L		1.44	4.8	8/09/2001	JF
SW-846-8270C	Dibenzo(a,h)anthracene	<1.5	ug/L		1.5	5	8/09/2001	JF
SW-846-8270C	Dibenzofuran	4.1	ug/L	13	1.28	4.2666	8/09/2001	JF
SW-846-8270C	Diethylphthalate	<1.3	ug/L		1.26	4.2	8/09/2001	JF
SW-846-8270C	Dimethylphthalate	<1.5	ug/L		1.52	5.0666	8/09/2001	JF
SW-846-8270C	Diphenylamine	<2.6	ug/L		2.6	8.6666	8/09/2001	JF
SW-846-8270C	Ethyl methanesulfonate	<1.7	ug/L		1.7	5.6666	8/09/2001	JF
SW-846-8270C	Fluoranthene	<2.0	ug/L		1.98	6.6	8/09/2001	JF
SW-846-8270C	Fluorene	6.3	ug/L		1.2	4	8/09/2001	JF
SW-846-8270C	Hexachlorobenzene	<1.6	ug/L		1.58	5.2666	8/09/2001	JF
SW-846-8270C	Hexachlorobutadiene	<7.6	ug/L		7.6	25.333	8/09/2001	JF
SW-846-8270C	Hexachlorocyclopentadiene	<5.4	ug/L		5.4	18	8/09/2001	JF
SW-846-8270C	Hexachloroethane	<5.2	ug/L		5.2	17.333	8/09/2001	JF
SW-846-8270C	Indeno(1,2,3-cd)pyrene	<1.4	ug/L		1.44	4.8	8/09/2001	JF
SW-846-8270C	Isophorone	<1.4	ug/L		1.42	4.7333	8/09/2001	JF
SW-846-8270C	Methyl methanesulfonate	<1.1	ug/L		1.06	3.5333	8/09/2001	JF
SW-846-8270C	n-Nitrosodi-n-butylamine	<1.5	ug/L		1.46	4.8666	8/09/2001	JF
SW-846-8270C	n-Nitrosodi-n-propylamine	<1.4	ug/L		1.4	4.6666	8/09/2001	JF
SW-846-8270C	n-Nitrosodimethylamine	<1.9	ug/L		1.88	6.2666	8/09/2001	JF
SW-846-8270C	n-Nitrosodiphenylamine	<2.6	ug/L		2.6	8.6666	8/09/2001	JF
SW-846-8270C	n-Nitrosopiperidine	<1.4	ug/L		1.4	4.6666	8/09/2001	JF
SW-846-8270C	Naphthalene	60	ug/L		1.38	4.6	8/09/2001	JF
SW-846-8270C	Nitrobenzene	<1.4	ug/L		1.42	4.7333	8/09/2001	JF
SW-846-8270C	Nitrobenzene-d5 - Surrogate	61	% Rec		0	0	8/09/2001	JF
SW-846-8270C	p-Dimethylaminoazobenzene	<2.0	ug/L		2	6.6666	8/09/2001	JF
SW-846-8270C	Pentachlorobenzene	<2.2	ug/L		2.2	7.3333	8/09/2001	JF
SW-846-8270C	Pentachloronitrobenzene	<1.5	ug/L		1.54	5.1333	8/09/2001	JF
SW-846-8270C	Pentachlorophenol	1190	ug/L		1.72	5.7333	8/09/2001	JF
SW-846-8270C	Phenacetin	<1.8	ug/L		1.78	5.9333	8/09/2001	JF
SW-846-8270C	Phenanthrene	1.6	ug/L	13	1.42	4.7333	8/09/2001	JF
SW-846-8270C	Phenol	<1.1	ug/L		1.1	3.6666	8/09/2001	JF
SW-846-8270C	Phenol-d5 - Surrogate	44	% Rec		0	0	8/09/2001	JF
SW-846-8270C	Pronamide	<1.7	ug/L		1.68	5.6	8/09/2001	JF
SW-846-8270C	Pyrene	<2.2	ug/L		2.2	7.3333	8/09/2001	JF
SW-846-8270C	Pyridine	<3.0	ug/L		3	10	8/09/2001	JF
SW-846-8270C	Terphenyl-d14 - Surrogate	56	% Rec		0	0	8/09/2001	JF
01REL010547	6/27/2001	MW-10						
EPA 1613	Dioxin Analysis	See Attached			0	0	7/22/2001	PJK
SW-846-8270C	Extraction Date	Complete			0	0	7/03/2001	JF
SW-846-8270C	1,2,4,5-Tetrachlorobenzene	<3.0	ug/L		3	10	8/09/2001	JF
SW-846-8270C	1,2,4-Trichlorobenzene	<3.6	ug/L		3.6	12	8/09/2001	JF
SW-846-8270C	1,2-Dichlorobenzene	<3.2	ug/L		3.2	10.666	8/09/2001	JF

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Chain: 85968

Report Date: 8/23/2001

Method	Parameter Name	Result	Units	Flag	MDL	PQL	Anl Date	Analyst
Lab No.	Collect Date	Sample ID						
SW-846-8270C	1,2-Diphenylhydrazine	<1.3	ug/L		1.34	4.4666	8/09/2001	JF
SW-846-8270C	1,3-Dichlorobenzene	<3.6	ug/L		3.6	12	8/09/2001	JF
SW-846-8270C	1,4-Dichlorobenzene	<3.4	ug/L		3.4	11.333	8/09/2001	JF
SW-846-8270C	1-Chloronaphthalene	<1.8	ug/L		1.84	6.1333	8/09/2001	JF
SW-846-8270C	1-Naphthylamine	<2.2	ug/L		2.2	7.3333	8/09/2001	JF
SW-846-8270C	2,3,4,6-Tetrachlorophenol	<1.4	ug/L		1.36	4.5333	8/09/2001	JF
SW-846-8270C	2,4,5-Trichlorophenol	<1.6	ug/L		1.64	5.4666	8/09/2001	JF
SW-846-8270C	2,4,6-Tribromophenol - Surrogate	77	% Rec		0	0	8/09/2001	JF
SW-846-8270C	2,4,6-Trichlorophenol	<1.6	ug/L		1.62	5.4	8/09/2001	JF
SW-846-8270C	2,4-Dichlorophenol	<1.5	ug/L		1.54	5.1333	8/09/2001	JF
SW-846-8270C	2,4-Dimethylphenol	<2.8	ug/L		2.8	9.3333	8/09/2001	JF
SW-846-8270C	2,4-Dinitrophenol	<1.8	ug/L		1.78	5.9333	8/09/2001	JF
SW-846-8270C	2,4-Dinitrotoluene	<1.6	ug/L		1.58	5.2666	8/09/2001	JF
SW-846-8270C	2,6-Dichlorophenol	<1.5	ug/L		1.52	5.0666	8/09/2001	JF
SW-846-8270C	2,6-Dinitrotoluene	<1.6	ug/L		1.6	5.3333	8/09/2001	JF
SW-846-8270C	2-Chloronaphthalene	<1.6	ug/L		1.56	5.2	8/09/2001	JF
SW-846-8270C	2-Chlorophenol	<1.2	ug/L		1.2	4	8/09/2001	JF
SW-846-8270C	2-Fluorobiphenyl - Surrogate	64	% Rec		0	0	8/09/2001	JF
SW-846-8270C	2-Fluorophenol - Surrogate	50	% Rec		0	0	8/09/2001	JF
SW-846-8270C	2-Methylnaphthalene	8.6	ug/L		1.6	5.3333	8/09/2001	JF
SW-846-8270C	2-Methylphenol	<1.7	ug/L		1.7	5.6666	8/09/2001	JF
SW-846-8270C	2-Naphthylamine	<2.2	ug/L		2.2	7.3333	8/09/2001	JF
SW-846-8270C	2-Nitroaniline	<1.6	ug/L		1.64	5.4666	8/09/2001	JF
SW-846-8270C	2-Nitrophenol	<1.3	ug/L		1.28	4.2666	8/09/2001	JF
SW-846-8270C	2-Picoline	<3.0	ug/L		3	10	8/09/2001	JF
SW-846-8270C	3 & 4-Methylphenol	<1.8	ug/L		1.76	5.8666	8/09/2001	JF
SW-846-8270C	3,3'-Dichlorobenzidine	<0.84	ug/L		0.84	2.8	8/09/2001	JF
SW-846-8270C	3-Methylcholanthrene	<1.9	ug/L		1.94	6.4666	8/09/2001	JF
SW-846-8270C	3-Nitroaniline	<2.2	ug/L		2.2	7.3333	8/09/2001	JF
SW-846-8270C	4,6-Dinitro-2-methylphenol	<2.6	ug/L		2.6	8.6666	8/09/2001	JF
SW-846-8270C	4-Aminobiphenyl	<1.5	ug/L		1.5	5	8/09/2001	JF
SW-846-8270C	4-Bromophenyl phenyl ether	<1.2	ug/L		1.2	4	8/09/2001	JF
SW-846-8270C	4-Chloro-3-methylphenol	<1.7	ug/L		1.7	5.6666	8/09/2001	JF
SW-846-8270C	4-Chloroaniline	<1.3	ug/L		1.32	4.4	8/09/2001	JF
SW-846-8270C	4-Chlorophenyl phenyl ether	<1.2	ug/L		1.24	4.1333	8/09/2001	JF
SW-846-8270C	4-Nitroaniline	<1.3	ug/L		1.32	4.4	8/09/2001	JF
SW-846-8270C	4-Nitrophenol	<1.9	ug/L		1.86	6.2	8/09/2001	JF
SW-846-8270C	7,12-Dimethylbenz(a)anthracene	<2.0	ug/L		2	6.6666	8/09/2001	JF
SW-846-8270C	Acenaphthene	<1.1	ug/L		1.12	3.7333	8/09/2001	JF
SW-846-8270C	Acenaphthylene	<1.5	ug/L		1.46	4.8666	8/09/2001	JF
SW-846-8270C	Acetophenone	<1.6	ug/L		1.6	5.3333	8/09/2001	JF
SW-846-8270C	Aniline	<1.7	ug/L		1.74	5.8	8/09/2001	JF
SW-846-8270C	Anthracene	<1.3	ug/L		1.3	4.3333	8/09/2001	JF
SW-846-8270C	Benzidine	<2.8	ug/L		2.8	9.3333	8/09/2001	JF
SW-846-8270C	Benzo(a)anthracene	<1.4	ug/L		1.42	4.7333	8/09/2001	JF
SW-846-8270C	Benzo(a)pyrene	<1.1	ug/L		1.1	3.6666	8/09/2001	JF

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 Client ID: L14
 Chain: 85968
 Report Date: 8/23/2001

Method	Parameter Name	Result	Units	Flag	MDL	PQL	Anal Date	Analyst	
Lab No	Collect Date	Sample ID							
SW-846-8270C		Benzo(b)fluoranthene	<1.3		ug/L	1.28	4.2666	8/09/2001	JF
SW-846-8270C		Benzo(g,h,i)perylene	<1.3		ug/L	1.32	4.4	8/09/2001	JF
SW-846-8270C		Benzo(k)fluoranthene	<1.6		ug/L	1.58	5.2666	8/09/2001	JF
SW-846-8270C		Benzyl alcohol	<1.8		ug/L	1.8	6	8/09/2001	JF
SW-846-8270C		bis(2-Chloroethoxy)methane	<1.5		ug/L	1.48	4.9333	8/09/2001	JF
SW-846-8270C		bis(2-Chloroethyl)ether	<1.9		ug/L	1.9	6.3333	8/09/2001	JF
SW-846-8270C		bis(2-Chloroisopropyl)ether	<1.5		ug/L	1.54	5.1333	8/09/2001	JF
SW-846-8270C		bis(2-Ethylhexyl)phthalate	<4.6		ug/L	4.6	15.333	8/09/2001	JF
SW-846-8270C		Butylbenzylphthalate	<1.8		ug/L	1.8	6	8/09/2001	JF
SW-846-8270C		Chrysene	<1.6		ug/L	1.56	5.2	8/09/2001	JF
SW-846-8270C		Di-n-butylphthalate	<1.5		ug/L	1.52	5.0666	8/09/2001	JF
SW-846-8270C		Di-n-octylphthalate	<1.5		ug/L	1.52	5.0666	8/09/2001	JF
SW-846-8270C		Dibenz(a,j)acridine	<1.4		ug/L	1.44	4.8	8/09/2001	JF
SW-846-8270C		Dibenzo(a,h)anthracene	<1.5		ug/L	1.5	5	8/09/2001	JF
SW-846-8270C		Dibenzofuran	<1.3		ug/L	1.28	4.2666	8/09/2001	JF
SW-846-8270C		Diethylphthalate	<1.3		ug/L	1.26	4.2	8/09/2001	JF
SW-846-8270C		Dimethylphthalate	<1.5		ug/L	1.52	5.0666	8/09/2001	JF
SW-846-8270C		Diphenylamine	<2.6		ug/L	2.6	8.6666	8/09/2001	JF
SW-846-8270C		Ethyl methanesulfonate	<1.7		ug/L	1.7	5.6666	8/09/2001	JF
SW-846-8270C		Fluoranthene	<2.0		ug/L	1.98	6.6	8/09/2001	JF
SW-846-8270C		Fluorene	<1.2		ug/L	1.2	4	8/09/2001	JF
SW-846-8270C		Hexachlorobenzene	<1.6		ug/L	1.58	5.2666	8/09/2001	JF
SW-846-8270C		Hexachlorobutadiene	<7.6		ug/L	7.6	25.333	8/09/2001	JF
SW-846-8270C		Hexachlorocyclopentadiene	<5.4		ug/L	5.4	18	8/09/2001	JF
SW-846-8270C		Hexachloroethane	<5.2		ug/L	5.2	17.333	8/09/2001	JF
SW-846-8270C		Indeno(1,2,3-cd)pyrene	<1.4		ug/L	1.44	4.8	8/09/2001	JF
SW-846-8270C		Isophorone	<1.4		ug/L	1.42	4.7333	8/09/2001	JF
SW-846-8270C		Methyl methanesulfonate	<1.1		ug/L	1.06	3.5333	8/09/2001	JF
SW-846-8270C		n-Nitrosodi-n-butylamine	<1.5		ug/L	1.46	4.8666	8/09/2001	JF
SW-846-8270C		n-Nitrosodi-n-propylamine	<1.4		ug/L	1.4	4.6666	8/09/2001	JF
SW-846-8270C		n-Nitrosodimethylamine	<1.9		ug/L	1.88	6.2666	8/09/2001	JF
SW-846-8270C		n-Nitrosodiphenylamine	<2.6		ug/L	2.6	8.6666	8/09/2001	JF
SW-846-8270C		n-Nitrosopiperidine	<1.4		ug/L	1.4	4.6666	8/09/2001	JF
SW-846-8270C		Naphthalene	<1.4		ug/L	1.38	4.6	8/09/2001	JF
SW-846-8270C		Nitrobenzene	<1.4		ug/L	1.42	4.7333	8/09/2001	JF
SW-846-8270C		Nitrobenzene-d5 - Surrogate	60		% Rec	0	0	8/09/2001	JF
SW-846-8270C		p-Dimethylaminoazobenzene	<2.0		ug/L	2	6.6666	8/09/2001	JF
SW-846-8270C		Pentachlorobenzene	<2.2		ug/L	2.2	7.3333	8/09/2001	JF
SW-846-8270C		Pentachloronitrobenzene	<1.5		ug/L	1.54	5.1333	8/09/2001	JF
SW-846-8270C		Pentachlorophenol	58		ug/L	1.72	5.7333	8/09/2001	JF
SW-846-8270C		Phenacetin	<1.8		ug/L	1.78	5.9333	8/09/2001	JF
SW-846-8270C		Phenanthrene	<1.4		ug/L	1.42	4.7333	8/09/2001	JF
SW-846-8270C		Phenol	<1.1		ug/L	1.1	3.6666	8/09/2001	JF
SW-846-8270C		Phenol-d5 - Surrogate	44		% Rec	0	0	8/09/2001	JF
SW-846-8270C		Pronamide	<1.7		ug/L	1.68	5.6	8/09/2001	JF
SW-846-8270C		Pyrene	<2.2		ug/L	2.2	7.3333	8/09/2001	JF

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Report Date: 8/23/2001

Method	Parameter Name	Result	Units	Flag	MDL	PQL	Anal Date	Analyst
Lab No	Collect Date	Sample ID						
SW-846-8270C	Pyridine	<3.0	ug/L		3	10	8/09/2001	JF
SW-846-8270C	Terphenyl-d14 - Surrogate	56	% Rec		0	0	8/09/2001	JF
01REL010548	6/27/2001	DMW-1						
EPA 1613	Dioxin Analysis	See Attached			0	0	7/22/2001	PJK
SW-846-8270C	Extraction Date	Complete			0	0	7/03/2001	JF
SW-846-8270C	1,2,4,5-Tetrachlorobenzene	<3.0	ug/L		3	10	8/10/2001	JF
SW-846-8270C	1,2,4-Trichlorobenzene	<3.6	ug/L		3.6	12	8/10/2001	JF
SW-846-8270C	1,2-Dichlorobenzene	<3.2	ug/L		3.2	10.666	8/10/2001	JF
SW-846-8270C	1,2-Diphenylhydrazine	<1.3	ug/L		1.34	4.4666	8/10/2001	JF
SW-846-8270C	1,3-Dichlorobenzene	<3.6	ug/L		3.6	12	8/10/2001	JF
SW-846-8270C	1,4-Dichlorobenzene	<3.4	ug/L		3.4	11.333	8/10/2001	JF
SW-846-8270C	1-Chloronaphthalene	<1.8	ug/L		1.84	6.1333	8/10/2001	JF
SW-846-8270C	1-Naphthylamine	<2.2	ug/L		2.2	7.3333	8/10/2001	JF
SW-846-8270C	2,3,4,6-Tetrachlorophenol	264	ug/L		1.36	4.5333	8/10/2001	JF
SW-846-8270C	2,4,5-Trichlorophenol	<1.6	ug/L		1.64	5.4666	8/10/2001	JF
SW-846-8270C	2,4,6-Tribromophenol - Surrogate	72	% Rec		0	0	8/10/2001	JF
SW-846-8270C	2,4,6-Trichlorophenol	<1.6	ug/L		1.62	5.4	8/10/2001	JF
SW-846-8270C	2,4-Dichlorophenol	<1.5	ug/L		1.54	5.1333	8/10/2001	JF
SW-846-8270C	2,4-Dimethylphenol	<2.8	ug/L		2.8	9.3333	8/10/2001	JF
SW-846-8270C	2,4-Dinitrophenol	<1.8	ug/L		1.78	5.9333	8/10/2001	JF
SW-846-8270C	2,4-Dinitrotoluene	<1.6	ug/L		1.58	5.2666	8/10/2001	JF
SW-846-8270C	2,6-Dichlorophenol	<1.5	ug/L		1.52	5.0666	8/10/2001	JF
SW-846-8270C	2,6-Dinitrotoluene	<1.6	ug/L		1.6	5.3333	8/10/2001	JF
SW-846-8270C	2-Chloronaphthalene	<1.6	ug/L		1.56	5.2	8/10/2001	JF
SW-846-8270C	2-Chlorophenol	<1.2	ug/L		1.2	4	8/10/2001	JF
SW-846-8270C	2-Fluorobiphenyl - Surrogate	68	% Rec		0	0	8/10/2001	JF
SW-846-8270C	2-Fluorophenol - Surrogate	49	% Rec		0	0	8/10/2001	JF
SW-846-8270C	2-Methylnaphthalene	37	ug/L		1.6	5.3333	8/10/2001	JF
SW-846-8270C	2-Methylphenol	<1.7	ug/L		1.7	5.6666	8/10/2001	JF
SW-846-8270C	2-Naphthylamine	<2.2	ug/L		2.2	7.3333	8/10/2001	JF
SW-846-8270C	2-Nitroaniline	<1.6	ug/L		1.64	5.4666	8/10/2001	JF
SW-846-8270C	2-Nitrophenol	<1.3	ug/L		1.28	4.2666	8/10/2001	JF
SW-846-8270C	2-Picoline	<3.0	ug/L		3	10	8/10/2001	JF
SW-846-8270C	3 & 4-Methylphenol	<1.8	ug/L		1.76	5.8666	8/10/2001	JF
SW-846-8270C	3,3'-Dichlorobenzidine	<0.84	ug/L		0.84	2.8	8/10/2001	JF
SW-846-8270C	3-Methylcholanthrene	<1.9	ug/L		1.94	6.4666	8/10/2001	JF
SW-846-8270C	3-Nitroaniline	<2.2	ug/L		2.2	7.3333	8/10/2001	JF
SW-846-8270C	4,6-Dinitro-2-methylphenol	<2.6	ug/L		2.6	8.6666	8/10/2001	JF
SW-846-8270C	4-Aminobiphenyl	<1.5	ug/L		1.5	5	8/10/2001	JF
SW-846-8270C	4-Bromophenyl phenyl ether	<1.2	ug/L		1.2	4	8/10/2001	JF
SW-846-8270C	4-Chloro-3-methylphenol	<1.7	ug/L		1.7	5.6666	8/10/2001	JF
SW-846-8270C	4-Chloroaniline	<1.3	ug/L		1.32	4.4	8/10/2001	JF
SW-846-8270C	4-Chlorophenyl phenyl ether	<1.2	ug/L		1.24	4.1333	8/10/2001	JF
SW-846-8270C	4-Nitroaniline	<1.3	ug/L		1.32	4.4	8/10/2001	JF
SW-846-8270C	4-Nitrophenol	<1.9	ug/L		1.86	6.2	8/10/2001	JF
SW-846-8270C	7,12-Dimethylbenz(a)anthracene	<2.0	ug/L		2	6.6666	8/10/2001	JF

Robert E. Lee & Associates, Inc.
Wisconsin Certification Number: 405043870
Certificate of Analysis Report

Robert E Lee & Associates, Inc
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Green Bay, WI 54301-2878
Project Number: 13551002
Project Name: WEISENBERGER TIE & LUMBER

Client ID: L14
Chain: 85968
Report Date: 8/23/2001

Method	Parameter Name	Result	Units	Flag	MDL	PQL	Anl Date	Analyst
Lab No.	Collect Date	Sample ID						
SW-846-8270C	Acenaphthene	24	ug/L		1.12	3.7333	8/10/2001	JF
SW-846-8270C	Acenaphthylene	<1.5	ug/L		1.46	4.8666	8/10/2001	JF
SW-846-8270C	Acetophenone	<1.6	ug/L		1.6	5.3333	8/10/2001	JF
SW-846-8270C	Aniline	<1.7	ug/L		1.74	5.8	8/10/2001	JF
SW-846-8270C	Anthracene	<1.3	ug/L		1.3	4.3333	8/10/2001	JF
SW-846-8270C	Benzidine	<2.8	ug/L		2.8	9.3333	8/10/2001	JF
SW-846-8270C	Benzo(a)anthracene	<1.4	ug/L		1.42	4.7333	8/10/2001	JF
SW-846-8270C	Benzo(a)pyrene	<1.1	ug/L		1.1	3.6666	8/10/2001	JF
SW-846-8270C	Benzo(b)fluoranthene	<1.3	ug/L		1.28	4.2666	8/10/2001	JF
SW-846-8270C	Benzo(g,h,i)perylene	<1.3	ug/L		1.32	4.4	8/10/2001	JF
SW-846-8270C	Benzo(k)fluoranthene	<1.6	ug/L		1.58	5.2666	8/10/2001	JF
SW-846-8270C	Benzyl alcohol	<1.8	ug/L		1.8	6	8/10/2001	JF
SW-846-8270C	bis(2-Chloroethoxy)methane	<1.5	ug/L		1.48	4.9333	8/10/2001	JF
SW-846-8270C	bis(2-Chloroethyl)ether	<1.9	ug/L		1.9	6.3333	8/10/2001	JF
SW-846-8270C	bis(2-Chloroisopropyl)ether	<1.5	ug/L		1.54	5.1333	8/10/2001	JF
SW-846-8270C	bis(2-Ethylhexyl)phtthalate	<4.6	ug/L		4.6	15.333	8/10/2001	JF
SW-846-8270C	Butylbenzylphtthalate	<1.8	ug/L		1.8	6	8/10/2001	JF
SW-846-8270C	Chrysene	<1.6	ug/L		1.56	5.2	8/10/2001	JF
SW-846-8270C	Di-n-butylphtthalate	<1.5	ug/L		1.52	5.0666	8/10/2001	JF
SW-846-8270C	Di-n-octylphtthalate	<1.5	ug/L		1.52	5.0666	8/10/2001	JF
SW-846-8270C	Dibenz(a,j)acridine	<1.4	ug/L		1.44	4.8	8/10/2001	JF
SW-846-8270C	Dibenzo(a,h)anthracene	<1.5	ug/L		1.5	5	8/10/2001	JF
SW-846-8270C	Dibenzofuran	11	ug/L		1.28	4.2666	8/10/2001	JF
SW-846-8270C	Diethylphtthalate	<1.3	ug/L		1.26	4.2	8/10/2001	JF
SW-846-8270C	Dimethylphtthalate	<1.5	ug/L		1.52	5.0666	8/10/2001	JF
SW-846-8270C	Diphenylamine	<2.6	ug/L		2.6	8.6666	8/10/2001	JF
SW-846-8270C	Ethyl methanesulfonate	<1.7	ug/L		1.7	5.6666	8/10/2001	JF
SW-846-8270C	Fluoranthene	4.2	ug/L	13	1.98	6.6	8/10/2001	JF
SW-846-8270C	Fluorene	24	ug/L		1.2	4	8/10/2001	JF
SW-846-8270C	Hexachlorobenzene	<1.6	ug/L		1.58	5.2666	8/10/2001	JF
SW-846-8270C	Hexachlorobutadiene	<7.6	ug/L		7.6	25.333	8/10/2001	JF
SW-846-8270C	Hexachlorocyclopentadiene	<5.4	ug/L		5.4	18	8/10/2001	JF
SW-846-8270C	Hexachloroethane	<5.2	ug/L		5.2	17.333	8/10/2001	JF
SW-846-8270C	Indeno(1,2,3-cd)pyrene	<1.4	ug/L		1.44	4.8	8/10/2001	JF
SW-846-8270C	Isophorone	<1.4	ug/L		1.42	4.7333	8/10/2001	JF
SW-846-8270C	Methyl methanesulfonate	<1.1	ug/L		1.06	3.5333	8/10/2001	JF
SW-846-8270C	n-Nitrosodi-n-butylamine	<1.5	ug/L		1.46	4.8666	8/10/2001	JF
SW-846-8270C	n-Nitrosodi-n-propylamine	<1.4	ug/L		1.4	4.6666	8/10/2001	JF
SW-846-8270C	n-Nitrosodimethylamine	<1.9	ug/L		1.88	6.2666	8/10/2001	JF
SW-846-8270C	n-Nitrosodiphenylamine	<2.6	ug/L		2.6	8.6666	8/10/2001	JF
SW-846-8270C	n-Nitrosopiperidine	<1.4	ug/L		1.4	4.6666	8/10/2001	JF
SW-846-8270C	Naphthalene	<1.4	ug/L		1.38	4.6	8/10/2001	JF
SW-846-8270C	Nitrobenzene	<1.4	ug/L		1.42	4.7333	8/10/2001	JF
SW-846-8270C	Nitrobenzene-d5 - Surrogate	60	% Rec		0	0	8/10/2001	JF
SW-846-8270C	p-Dimethylaminoazobenzene	<2.0	ug/L		2	6.6666	8/10/2001	JF
SW-846-8270C	Pentachlorobenzene	<2.2	ug/L		2.2	7.3333	8/10/2001	JF

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Client ID: L14

Chain: 85968

Report Date: 8/23/2001

Method	Parameter Name	Result	Units	Flag	MDL	PQL	Anl Date	Analyst
Lab No.	Collect Date	Sample ID						
SW-846-8270C	Pentachloronitrobenzene	<1.5	ug/L		1.54	5.1333	8/10/2001	JF
SW-846-8270C	Pentachlorophenol	4050	ug/L	<u>21</u>	1.72	5.7333	8/10/2001	JF
SW-846-8270C	Phenacetin	<1.8	ug/L		1.78	5.9333	8/10/2001	JF
SW-846-8270C	Phenanthrene	29	ug/L		1.42	4.7333	8/10/2001	JF
SW-846-8270C	Phenol	<1.1	ug/L		1.1	3.6666	8/10/2001	JF
SW-846-8270C	Phenol-d5 - Surrogate	46	% Rec		0	0	8/10/2001	JF
SW-846-8270C	Pronamide	<1.7	ug/L		1.68	5.6	8/10/2001	JF
SW-846-8270C	Pyrene	5.4	ug/L	<u>13</u>	2.2	7.3333	8/10/2001	JF
SW-846-8270C	Pyridine	<3.0	ug/L		3	10	8/10/2001	JF
SW-846-8270C	Terphenyl-d14 - Surrogate	52	% Rec		0	0	8/10/2001	JF
01REL010549	6/27/2001	DMW-2						
EPA 1613	Dioxin Analysis	See Attached			0	0	7/22/2003	PJK
SW-846-8270C	Extraction Date	Complete			0	0	7/03/2001	JF
SW-846-8270C	1,2,4,5-Tetrachlorobenzene	<3.0	ug/L		3	10	8/10/2001	JF
SW-846-8270C	1,2,4-Trichlorobenzene	<3.6	ug/L		3.6	12	8/10/2001	JF
SW-846-8270C	1,2-Dichlorobenzene	<3.2	ug/L		3.2	10.666	8/10/2001	JF
SW-846-8270C	1,2-Diphenylhydrazine	<1.3	ug/L		1.34	4.4666	8/10/2001	JF
SW-846-8270C	1,3-Dichlorobenzene	<3.6	ug/L		3.6	12	8/10/2001	JF
SW-846-8270C	1,4-Dichlorobenzene	<3.4	ug/L		3.4	11.333	8/10/2001	JF
SW-846-8270C	1-Chloronaphthalene	<1.8	ug/L		1.84	6.1333	8/10/2001	JF
SW-846-8270C	1-Naphthylamine	<2.2	ug/L		2.2	7.3333	8/10/2001	JF
SW-846-8270C	2,3,4,6-Tetrachlorophenol	<1.4	ug/L		1.36	4.5333	8/10/2001	JF
SW-846-8270C	2,4,5-Trichlorophenol	<1.6	ug/L		1.64	5.4666	8/10/2001	JF
SW-846-8270C	2,4,6-Tribromophenol - Surrogate	64	% Rec		0	0	8/10/2001	JF
SW-846-8270C	2,4,6-Trichlorophenol	<1.6	ug/L		1.62	5.4	8/10/2001	JF
SW-846-8270C	2,4-Dichlorophenol	<1.5	ug/L		1.54	5.1333	8/10/2001	JF
SW-846-8270C	2,4-Dimethylphenol	<2.8	ug/L		2.8	9.3333	8/10/2001	JF
SW-846-8270C	2,4-Dinitrophenol	<1.8	ug/L		1.78	5.9333	8/10/2001	JF
SW-846-8270C	2,4-Dinitrotoluene	<1.6	ug/L		1.58	5.2666	8/10/2001	JF
SW-846-8270C	2,6-Dichlorophenol	<1.5	ug/L		1.52	5.0666	8/10/2001	JF
SW-846-8270C	2,6-Dinitrotoluene	<1.6	ug/L		1.6	5.3333	8/10/2001	JF
SW-846-8270C	2-Chloronaphthalene	<1.6	ug/L		1.56	5.2	8/10/2001	JF
SW-846-8270C	2-Chlorophenol	<1.2	ug/L		1.2	4	8/10/2001	JF
SW-846-8270C	2-Fluorobiphenyl - Surrogate	61	% Rec		0	0	8/10/2001	JF
SW-846-8270C	2-Fluorophenol - Surrogate	47	% Rec		0	0	8/10/2001	JF
SW-846-8270C	2-Methylnaphthalene	<1.6	ug/L		1.6	5.3333	8/10/2001	JF
SW-846-8270C	2-Methylphenol	<1.7	ug/L		1.7	5.6666	8/10/2001	JF
SW-846-8270C	2-Naphthylamine	<2.2	ug/L		2.2	7.3333	8/10/2001	JF
SW-846-8270C	2-Nitroaniline	<1.6	ug/L		1.64	5.4666	8/10/2001	JF
SW-846-8270C	2-Nitrophenol	<1.3	ug/L		1.28	4.2666	8/10/2001	JF
SW-846-8270C	2-Picoline	<3.0	ug/L		3	10	8/10/2001	JF
SW-846-8270C	3 & 4-Methylphenol	<1.8	ug/L		1.76	5.8666	8/10/2001	JF
SW-846-8270C	3,3'-Dichlorobenzidine	<0.84	ug/L		0.84	2.8	8/10/2001	JF
SW-846-8270C	3-Methylcholanthrene	<1.9	ug/L		1.94	6.4666	8/10/2001	JF
SW-846-8270C	3-Nitroaniline	<2.2	ug/L		2.2	7.3333	8/10/2001	JF
SW-846-8270C	4,6-Dinitro-2-methylphenol	<2.6	ug/L		2.6	8.6666	8/10/2001	JF

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Client ID: L14

Chain: 85968

Report Date: 8/23/2001

Method	Parameter Name	Result	Units	Flag	MDL	PQL	Anl Date	Analyst
Lab No	Collect Date	Sample ID						
SW-846-8270C	4-Aminobiphenyl	<1.5	ug/L		1.5	5	8/10/2001	JF
SW-846-8270C	4-Bromophenyl phenyl ether	<1.2	ug/L		1.2	4	8/10/2001	JF
SW-846-8270C	4-Chloro-3-methylphenol	<1.7	ug/L		1.7	5.6666	8/10/2001	JF
SW-846-8270C	4-Chloroaniline	<1.3	ug/L		1.32	4.4	8/10/2001	JF
SW-846-8270C	4-Chlorophenyl phenyl ether	<1.2	ug/L		1.24	4.1333	8/10/2001	JF
SW-846-8270C	4-Nitroaniline	<1.3	ug/L		1.32	4.4	8/10/2001	JF
SW-846-8270C	4-Nitrophenol	<1.9	ug/L		1.86	6.2	8/10/2001	JF
SW-846-8270C	7,12-Dimethylbenz(a)anthracene	<2.0	ug/L		2	6.6666	8/10/2001	JF
SW-846-8270C	Acenaphthene	<1.1	ug/L		1.12	3.7333	8/10/2001	JF
SW-846-8270C	Acenaphthylene	<1.5	ug/L		1.46	4.8666	8/10/2001	JF
SW-846-8270C	Acetophenone	<1.6	ug/L		1.6	5.3333	8/10/2001	JF
SW-846-8270C	Aniline	<1.7	ug/L		1.74	5.8	8/10/2001	JF
SW-846-8270C	Anthracene	<1.3	ug/L		1.3	4.3333	8/10/2001	JF
SW-846-8270C	Benzidine	<2.8	ug/L		2.8	9.3333	8/10/2001	JF
SW-846-8270C	Benzo(a)anthracene	<1.4	ug/L		1.42	4.7333	8/10/2001	JF
SW-846-8270C	Benzo(a)pyrene	<1.1	ug/L		1.1	3.6666	8/10/2001	JF
SW-846-8270C	Benzo(b)fluoranthene	<1.3	ug/L		1.28	4.2666	8/10/2001	JF
SW-846-8270C	Benzo(g,h,i)perylene	<1.3	ug/L		1.32	4.4	8/10/2001	JF
SW-846-8270C	Benzo(k)fluoranthene	<1.6	ug/L		1.58	5.2666	8/10/2001	JF
SW-846-8270C	Benzyl alcohol	<1.8	ug/L		1.8	6	8/10/2001	JF
SW-846-8270C	bis(2-Chloroethoxy)methane	<1.5	ug/L		1.48	4.9333	8/10/2001	JF
SW-846-8270C	bis(2-Chloroethyl)ether	<1.9	ug/L		1.9	6.3333	8/10/2001	JF
SW-846-8270C	bis(2-Chloroisopropyl)ether	<1.5	ug/L		1.54	5.1333	8/10/2001	JF
SW-846-8270C	bis(2-Ethylhexyl)phthalate	<4.6	ug/L		4.6	15.333	8/10/2001	JF
SW-846-8270C	Butylbenzylphthalate	<1.8	ug/L		1.8	6	8/10/2001	JF
SW-846-8270C	Chrysene	<1.6	ug/L		1.56	5.2	8/10/2001	JF
SW-846-8270C	Di-n-butylphthalate	<1.5	ug/L		1.52	5.0666	8/10/2001	JF
SW-846-8270C	Di-n-octylphthalate	<1.5	ug/L		1.52	5.0666	8/10/2001	JF
SW-846-8270C	Dibenz(a,j)acridine	<1.4	ug/L		1.44	4.8	8/10/2001	JF
SW-846-8270C	Dibenzo(a,h)anthracene	<1.5	ug/L		1.5	5	8/10/2001	JF
SW-846-8270C	Dibenzofuran	<1.3	ug/L		1.28	4.2666	8/10/2001	JF
SW-846-8270C	Diethylphthalate	<1.3	ug/L		1.26	4.2	8/10/2001	JF
SW-846-8270C	Dimethylphthalate	<1.5	ug/L		1.52	5.0666	8/10/2001	JF
SW-846-8270C	Diphenylamine	<2.6	ug/L		2.6	8.6666	8/10/2001	JF
SW-846-8270C	Ethyl methanesulfonate	<1.7	ug/L		1.7	5.6666	8/10/2001	JF
SW-846-8270C	Fluoranthene	<2.0	ug/L		1.98	6.6	8/10/2001	JF
SW-846-8270C	Fluorene	<1.2	ug/L		1.2	4	8/10/2001	JF
SW-846-8270C	Hexachlorobenzene	<1.6	ug/L		1.58	5.2666	8/10/2001	JF
SW-846-8270C	Hexachlorobutadiene	<7.6	ug/L		7.6	25.333	8/10/2001	JF
SW-846-8270C	Hexachlorocyclopentadiene	<5.4	ug/L		5.4	18	8/10/2001	JF
SW-846-8270C	Hexachloroethane	<5.2	ug/L		5.2	17.333	8/10/2001	JF
SW-846-8270C	Indeno(1,2,3-cd)pyrene	<1.4	ug/L		1.44	4.8	8/10/2001	JF
SW-846-8270C	Isophorone	<1.4	ug/L		1.42	4.7333	8/10/2001	JF
SW-846-8270C	Methyl methanesulfonate	<1.1	ug/L		1.06	3.5333	8/10/2001	JF
SW-846-8270C	n-Nitrosodi-n-butylamine	<1.5	ug/L		1.46	4.8666	8/10/2001	JF
SW-846-8270C	n-Nitrosodi-n-propylamine	<1.4	ug/L		1.4	4.6666	8/10/2001	JF

Robert E. Lee & Associates, Inc.
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Robert E Lee & Associates, Inc
 2825 S Webster Ave

Attn: Jim Caine
 Phone: (920)336-6338
 Fax: (920)336-9141

Green Bay, WI 54301-2878
 Project Number: 13551002
 Project Name: WEISENBERGER TIE & LUMBER

Client ID: L14
 Chain: 85968
 Report Date: 8/23/2001

Method	Parameter Name	Result	Units	Flag	MDL	PQL	Anl Date	Analyst
Lab No	Collect Date	Sample ID						
SW-846-8270C	n-Nitrosodimethylamine	<1.9	ug/L		1.88	6.2666	8/10/2001	JF
SW-846-8270C	n-Nitrosodiphenylamine	<2.6	ug/L		2.6	8.6666	8/10/2001	JF
SW-846-8270C	n-Nitrosopiperidine	<1.4	ug/L		1.4	4.6666	8/10/2001	JF
SW-846-8270C	Naphthalene	<1.4	ug/L		1.38	4.6	8/10/2001	JF
SW-846-8270C	Nitrobenzene	<1.4	ug/L		1.42	4.7333	8/10/2001	JF
SW-846-8270C	Nitrobenzene-d5 - Surrogate	59	% Rec		0	0	8/10/2001	JF
SW-846-8270C	p-Dimethylaminoazobenzene	<2.0	ug/L		2	6.6666	8/10/2001	JF
SW-846-8270C	Pentachlorobenzene	<2.2	ug/L		2.2	7.3333	8/10/2001	JF
SW-846-8270C	Pentachloronitrobenzene	<1.5	ug/L		1.54	5.1333	8/10/2001	JF
SW-846-8270C	Pentachlorophenol	<1.7	ug/L		1.72	5.7333	8/10/2001	JF
SW-846-8270C	Phenacetin	<1.8	ug/L		1.78	5.9333	8/10/2001	JF
SW-846-8270C	Phenanthrene	<1.4	ug/L		1.42	4.7333	8/10/2001	JF
SW-846-8270C	Phenol	<1.1	ug/L		1.1	3.6666	8/10/2001	JF
SW-846-8270C	Phenol-d5 - Surrogate	42	% Rec		0	0	8/10/2001	JF
SW-846-8270C	Pronamide	<1.7	ug/L		1.68	5.6	8/10/2001	JF
SW-846-8270C	Pyrene	<2.2	ug/L		2.2	7.3333	8/10/2001	JF
SW-846-8270C	Pyridine	<3.0	ug/L		3	10	8/10/2001	JF
SW-846-8270C	Terphenyl-d14 - Surrogate	58	% Rec		0	0	8/10/2001	JF

01REL010550 6/27/2001 DMW-4

EPA 1613	Dioxin Analysis	See Attached			0	0	7/24/2001	PJK
SW-846-8021B	1,2,4-Trimethylbenzene	2.9	ug/L		0.23	0.7666	7/10/2001	CRW
SW-846-8021B	1,3,5-Trimethylbenzene	1.5	ug/L		0.21	0.7	7/10/2001	CRW
SW-846-8021B	Benzene	<0.21	ug/L		0.21	0.7	7/10/2001	CRW
SW-846-8021B	Ethylbenzene	<0.23	ug/L		0.23	0.7666	7/10/2001	CRW
SW-846-8021B	Fluorobenzene-Surrogate	101	% Rec		0	0	7/10/2001	CRW
SW-846-8021B	Methyl-tertiary-butyl ether	<0.091	ug/L		0.091	0.3033	7/10/2001	CRW
SW-846-8021B	Toluene	<0.22	ug/L		0.22	0.7333	7/10/2001	CRW
SW-846-8021B	Xylenes-Total	2.9	ug/L		0.44	1.4666	7/10/2001	CRW
SW-846-8270C	Extraction Date	Complete			0	0	7/03/2001	JF
SW-846-8270C	1,2,4,5-Tetrachlorobenzene	<3.0	ug/L		3	10	7/30/2001	JF
SW-846-8270C	1,2,4-Trichlorobenzene	<3.6	ug/L		3.6	12	7/30/2001	JF
SW-846-8270C	1,2-Dichlorobenzene	<3.2	ug/L		3.2	10.666	7/30/2001	JF
SW-846-8270C	1,2-Diphenylhydrazine	<1.3	ug/L		1.34	4.4666	7/30/2001	JF
SW-846-8270C	1,3-Dichlorobenzene	<3.6	ug/L		3.6	12	7/30/2001	JF
SW-846-8270C	1,4-Dichlorobenzene	<3.4	ug/L		3.4	11.333	7/30/2001	JF
SW-846-8270C	1-Chloronaphthalene	<1.8	ug/L		1.84	6.1333	7/30/2001	JF
SW-846-8270C	1-Naphthylamine	<2.2	ug/L		2.2	7.3333	7/30/2001	JF
SW-846-8270C	2,3,4,6-Tetrachlorophenol	1.9	ug/L	13	1.36	4.5333	7/30/2001	JF
SW-846-8270C	2,4,5-Trichlorophenol	<1.6	ug/L		1.64	5.4666	7/30/2001	JF
SW-846-8270C	2,4,6-Tribromophenol - Surrogate	64	% Rec		0	0	7/30/2001	JF
SW-846-8270C	2,4,6-Trichlorophenol	<1.6	ug/L		1.62	5.4	7/30/2001	JF
SW-846-8270C	2,4-Dichlorophenol	<1.5	ug/L		1.54	5.1333	7/30/2001	JF
SW-846-8270C	2,4-Dimethylphenol	<2.8	ug/L		2.8	9.3333	7/30/2001	JF
SW-846-8270C	2,4-Dinitrophenol	<1.8	ug/L		1.78	5.9333	7/30/2001	JF
SW-846-8270C	2,4-Dinitrotoluene	<1.6	ug/L		1.58	5.2666	7/30/2001	JF
SW-846-8270C	2,6-Dichlorophenol	<1.5	ug/L		1.52	5.0666	7/30/2001	JF

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Project Number: 13551002

Project Name: WEISENBERGER TIE & LUMBER

Attn: Jim Caine

Phone: (920)336-6338

Fax: (920)336-9141

Client ID: L14

Chain: 85968

Report Date: 8/23/2001

Method	Parameter Name	Result	Units	Flag	MDL	PQL	AnalDate	Analyst
Lab No.	Collect Date	Sample ID						
SW-846-8270C	2,6-Dinitrotoluene	<1.6	ug/L		1.6	5.3333	7/30/2001	JF
SW-846-8270C	2-Chloronaphthalene	<1.6	ug/L		1.56	5.2	7/30/2001	JF
SW-846-8270C	2-Chlorophenol	<1.2	ug/L		1.2	4	7/30/2001	JF
SW-846-8270C	2-Fluorobiphenyl - Surrogate	55	% Rec		0	0	7/30/2001	JF
SW-846-8270C	2-Fluorophenol - Surrogate	40	% Rec		0	0	7/30/2001	JF
SW-846-8270C	2-Methylnaphthalene	<1.6	ug/L		1.6	5.3333	7/30/2001	JF
SW-846-8270C	2-Methylphenol	<1.7	ug/L		1.7	5.6666	7/30/2001	JF
SW-846-8270C	2-Naphthylamine	<2.2	ug/L		2.2	7.3333	7/30/2001	JF
SW-846-8270C	2-Nitroaniline	<1.6	ug/L		1.64	5.4666	7/30/2001	JF
SW-846-8270C	2-Nitrophenol	<1.3	ug/L		1.28	4.2666	7/30/2001	JF
SW-846-8270C	2-Picoline	<3.0	ug/L		3	10	7/30/2001	JF
SW-846-8270C	3 & 4-Methylphenol	<1.8	ug/L		1.76	5.8666	7/30/2001	JF
SW-846-8270C	3,3'-Dichlorobenzidine	<0.84	ug/L		0.84	2.8	7/30/2001	JF
SW-846-8270C	3-Methylcholanthrene	<1.9	ug/L		1.94	6.4666	7/30/2001	JF
SW-846-8270C	3-Nitroaniline	<2.2	ug/L		2.2	7.3333	7/30/2001	JF
SW-846-8270C	4,6-Dinitro-2-methylphenol	<2.6	ug/L		2.6	8.6666	7/30/2001	JF
SW-846-8270C	4-Aminobiphenyl	<1.5	ug/L		1.5	5	7/30/2001	JF
SW-846-8270C	4-Bromophenyl phenyl ether	<1.2	ug/L		1.2	4	7/30/2001	JF
SW-846-8270C	4-Chloro-3-methylphenol	<1.7	ug/L		1.7	5.6666	7/30/2001	JF
SW-846-8270C	4-Chloroaniline	<1.3	ug/L		1.32	4.4	7/30/2001	JF
SW-846-8270C	4-Chlorophenyl phenyl ether	<1.2	ug/L		1.24	4.1333	7/30/2001	JF
SW-846-8270C	4-Nitroaniline	<1.3	ug/L		1.32	4.4	7/30/2001	JF
SW-846-8270C	4-Nitrophenol	<1.9	ug/L		1.86	6.2	7/30/2001	JF
SW-846-8270C	7,12-Dimethylbenz(a)anthracene	<2.0	ug/L		2	6.6666	7/30/2001	JF
SW-846-8270C	Acenaphthene	<1.1	ug/L		1.12	3.7333	7/30/2001	JF
SW-846-8270C	Acenaphthylene	<1.5	ug/L		1.46	4.8666	7/30/2001	JF
SW-846-8270C	Acetophenone	<1.6	ug/L		1.6	5.3333	7/30/2001	JF
SW-846-8270C	Aniline	<1.7	ug/L		1.74	5.8	7/30/2001	JF
SW-846-8270C	Anthracene	<1.3	ug/L		1.3	4.3333	7/30/2001	JF
SW-846-8270C	Benzidine	<2.8	ug/L		2.8	9.3333	7/30/2001	JF
SW-846-8270C	Benzo(a)anthracene	<1.4	ug/L		1.42	4.7333	7/30/2001	JF
SW-846-8270C	Benzo(a)pyrene	<1.1	ug/L		1.1	3.6666	7/30/2001	JF
SW-846-8270C	Benzo(b)fluoranthene	<1.3	ug/L		1.28	4.2666	7/30/2001	JF
SW-846-8270C	Benzo(g,h,i)perylene	<1.3	ug/L		1.32	4.4	7/30/2001	JF
SW-846-8270C	Benzo(k)fluoranthene	<1.6	ug/L		1.58	5.2666	7/30/2001	JF
SW-846-8270C	Benzyl alcohol	<1.8	ug/L		1.8	6	7/30/2001	JF
SW-846-8270C	bis(2-Chloroethoxy)methane	<1.5	ug/L		1.48	4.9333	7/30/2001	JF
SW-846-8270C	bis(2-Chloroethyl)ether	<1.9	ug/L		1.9	6.3333	7/30/2001	JF
SW-846-8270C	bis(2-Chloroisopropyl)ether	<1.5	ug/L		1.54	5.1333	7/30/2001	JF
SW-846-8270C	bis(2-Ethylhexyl)phthalate	<4.6	ug/L		4.6	15.333	7/30/2001	JF
SW-846-8270C	Butylbenzylphthalate	<1.8	ug/L		1.8	6	7/30/2001	JF
SW-846-8270C	Chrysene	<1.6	ug/L		1.56	5.2	7/30/2001	JF
SW-846-8270C	Di-n-butylphthalate	<1.5	ug/L		1.52	5.0666	7/30/2001	JF
SW-846-8270C	Di-n-octylphthalate	<1.5	ug/L		1.52	5.0666	7/30/2001	JF
SW-846-8270C	Dibenz(a,j)acridine	<1.4	ug/L		1.44	4.8	7/30/2001	JF
SW-846-8270C	Dibenzo(a,h)anthracene	<1.5	ug/L		1.5	5	7/30/2001	JF

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Green Bay, WI 54301-2878
 Project Number: 13551002
 Project Name: WEISENBERGER TIE & LUMBER

Method	Parameter Name	Result	Units	Flag	MDL	PQL	Anl Date	Analyst
Lab No.	Collect Date	Sample ID						
SW-846-8270C	Dibenzofuran	<1.3	ug/L		1.28	4.2666	7/30/2001	JF
SW-846-8270C	Diethylphthalate	<1.3	ug/L		1.26	4.2	7/30/2001	JF
SW-846-8270C	Dimethylphthalate	<1.5	ug/L		1.52	5.0666	7/30/2001	JF
SW-846-8270C	Diphenylamine	<2.6	ug/L		2.6	8.6666	7/30/2001	JF
SW-846-8270C	Ethyl methanesulfonate	<1.7	ug/L		1.7	5.6666	7/30/2001	JF
SW-846-8270C	Fluoranthene	<2.0	ug/L		1.98	6.6	7/30/2001	JF
SW-846-8270C	Fluorene	<1.2	ug/L		1.2	4	7/30/2001	JF
SW-846-8270C	Hexachlorobenzene	<1.6	ug/L		1.58	5.2666	7/30/2001	JF
SW-846-8270C	Hexachlorobutadiene	<7.6	ug/L		7.6	25.333	7/30/2001	JF
SW-846-8270C	Hexachlorocyclopentadiene	<5.4	ug/L		5.4	18	7/30/2001	JF
SW-846-8270C	Hexachloroethane	<5.2	ug/L		5.2	17.333	7/30/2001	JF
SW-846-8270C	Indeno(1,2,3-cd)pyrene	<1.4	ug/L		1.44	4.8	7/30/2001	JF
SW-846-8270C	Isophorone	<1.4	ug/L		1.42	4.7333	7/30/2001	JF
SW-846-8270C	Methyl methanesulfonate	<1.1	ug/L		1.06	3.5333	7/30/2001	JF
SW-846-8270C	n-Nitrosodi-n-butylamine	<1.5	ug/L		1.46	4.8666	7/30/2001	JF
SW-846-8270C	n-Nitrosodi-n-propylamine	<1.4	ug/L		1.4	4.6666	7/30/2001	JF
SW-846-8270C	n-Nitrosodimethylamine	<1.9	ug/L		1.88	6.2666	7/30/2001	JF
SW-846-8270C	n-Nitrosodiphenylamine	<2.6	ug/L		2.6	8.6666	7/30/2001	JF
SW-846-8270C	n-Nitrosopiperidine	<1.4	ug/L		1.4	4.6666	7/30/2001	JF
SW-846-8270C	Naphthalene	<1.4	ug/L		1.38	4.6	7/30/2001	JF
SW-846-8270C	Nitrobenzene	<1.4	ug/L		1.42	4.7333	7/30/2001	JF
SW-846-8270C	Nitrobenzene-d5 - Surrogate	47	% Rec		0	0	7/30/2001	JF
SW-846-8270C	p-Dimethylaminoazobenzene	<2.0	ug/L		2	6.6666	7/30/2001	JF
SW-846-8270C	Pentachlorobenzene	<2.2	ug/L		2.2	7.3333	7/30/2001	JF
SW-846-8270C	Pentachloronitrobenzene	<1.5	ug/L		1.54	5.1333	7/30/2001	JF
SW-846-8270C	Pentachlorophenol	38	ug/L		1.72	5.7333	7/30/2001	JF
SW-846-8270C	Phenacetin	<1.8	ug/L		1.78	5.9333	7/30/2001	JF
SW-846-8270C	Phenanthrene	<1.4	ug/L		1.42	4.7333	7/30/2001	JF
SW-846-8270C	Phenol	<1.1	ug/L		1.1	3.6666	7/30/2001	JF
SW-846-8270C	Phenol-d5 - Surrogate	37	% Rec		0	0	7/30/2001	JF
SW-846-8270C	Pronamide	<1.7	ug/L		1.68	5.6	7/30/2001	JF
SW-846-8270C	Pyrene	<2.2	ug/L		2.2	7.3333	7/30/2001	JF
SW-846-8270C	Pyridine	<3.0	ug/L		3	10	7/30/2001	JF
SW-846-8270C	Terphenyl-d14 - Surrogate	51	% Rec		0	0	7/30/2001	JF

01REL010551 6/27/2001 DMW-5

EPA 1613	Dioxin Analysis	See Attached			0	0	7/24/2001	PJK
SW-846-8021B	1,2,4-Trimethylbenzene	<0.23	ug/L		0.23	0.7666	7/10/2001	CRW
SW-846-8021B	1,3,5-Trimethylbenzene	<0.21	ug/L		0.21	0.7	7/10/2001	CRW
SW-846-8021B	Benzene	<0.21	ug/L		0.21	0.7	7/10/2001	CRW
SW-846-8021B	Ethylbenzene	<0.23	ug/L		0.23	0.7666	7/10/2001	CRW
SW-846-8021B	Fluorobenzene-Surrogate	101	% Rec		0	0	7/10/2001	CRW
SW-846-8021B	Methyl-tertiary-butyl ether	<0.091	ug/L		0.091	0.3033	7/10/2001	CRW
SW-846-8021B	Toluene	<0.22	ug/L		0.22	0.7333	7/10/2001	CRW
SW-846-8021B	Xylenes-Total	<0.44	ug/L		0.44	1.4666	7/10/2001	CRW
SW-846-8270C	Extraction Date	Complete			0	0	7/03/2001	JF
SW-846-8270C	1,2,4,5-Tetrachlorobenzene	<3.0	ug/L		3	10	7/30/2001	JF

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Method	Parameter Name	Result	Units	Flag	MDL	PQL	Anl Date	Analyst
Lab No	Collect Date	Sample ID						
SW-846-8270C	1,2,4-Trichlorobenzene	<3.6	ug/L		3.6	12	7/30/2001	JF
SW-846-8270C	1,2-Dichlorobenzene	<3.2	ug/L		3.2	10.666	7/30/2001	JF
SW-846-8270C	1,2-Diphenylhydrazine	<1.3	ug/L		1.34	4.4666	7/30/2001	JF
SW-846-8270C	1,3-Dichlorobenzene	<3.6	ug/L		3.6	12	7/30/2001	JF
SW-846-8270C	1,4-Dichlorobenzene	<3.4	ug/L		3.4	11.333	7/30/2001	JF
SW-846-8270C	1-Chloronaphthalene	<1.8	ug/L		1.84	6.1333	7/30/2001	JF
SW-846-8270C	1-Naphthylamine	<2.2	ug/L		2.2	7.3333	7/30/2001	JF
SW-846-8270C	2,3,4,6-Tetrachlorophenol	<1.4	ug/L		1.36	4.5333	7/30/2001	JF
SW-846-8270C	2,4,5-Trichlorophenol	<1.6	ug/L		1.64	5.4666	7/30/2001	JF
SW-846-8270C	2,4,6-Tribromophenol - Surrogate	68	% Rec		0	0	7/30/2001	JF
SW-846-8270C	2,4,6-Trichlorophenol	<1.6	ug/L		1.62	5.4	7/30/2001	JF
SW-846-8270C	2,4-Dichlorophenol	<1.5	ug/L		1.54	5.1333	7/30/2001	JF
SW-846-8270C	2,4-Dimethylphenol	<2.8	ug/L		2.8	9.3333	7/30/2001	JF
SW-846-8270C	2,4-Dinitrophenol	<1.8	ug/L		1.78	5.9333	7/30/2001	JF
SW-846-8270C	2,4-Dinitrotoluene	<1.6	ug/L		1.58	5.2666	7/30/2001	JF
SW-846-8270C	2,6-Dichlorophenol	<1.5	ug/L		1.52	5.0666	7/30/2001	JF
SW-846-8270C	2,6-Dinitrotoluene	<1.6	ug/L		1.6	5.3333	7/30/2001	JF
SW-846-8270C	2-Chloronaphthalene	<1.6	ug/L		1.56	5.2	7/30/2001	JF
SW-846-8270C	2-Chlorophenol	<1.2	ug/L		1.2	4	7/30/2001	JF
SW-846-8270C	2-Fluorobiphenyl - Surrogate	62	% Rec		0	0	7/30/2001	JF
SW-846-8270C	2-Fluorophenol - Surrogate	42	% Rec		0	0	7/30/2001	JF
SW-846-8270C	2-Methylnaphthalene	<1.6	ug/L		1.6	5.3333	7/30/2001	JF
SW-846-8270C	2-Methylphenol	<1.7	ug/L		1.7	5.6666	7/30/2001	JF
SW-846-8270C	2-Naphthylamine	<2.2	ug/L		2.2	7.3333	7/30/2001	JF
SW-846-8270C	2-Nitroaniline	<1.6	ug/L		1.64	5.4666	7/30/2001	JF
SW-846-8270C	2-Nitrophenol	<1.3	ug/L		1.28	4.2666	7/30/2001	JF
SW-846-8270C	2-Picoline	<3.0	ug/L		3	10	7/30/2001	JF
SW-846-8270C	3 & 4-Methylphenol	<1.8	ug/L		1.76	5.8666	7/30/2001	JF
SW-846-8270C	3,3'-Dichlorobenzidine	<0.84	ug/L		0.84	2.8	7/30/2001	JF
SW-846-8270C	3-Methylcholanthrene	<1.9	ug/L		1.94	6.4666	7/30/2001	JF
SW-846-8270C	3-Nitroaniline	<2.2	ug/L		2.2	7.3333	7/30/2001	JF
SW-846-8270C	4,6-Dinitro-2-methylphenol	<2.6	ug/L		2.6	8.6666	7/30/2001	JF
SW-846-8270C	4-Aminobiphenyl	<1.5	ug/L		1.5	5	7/30/2001	JF
SW-846-8270C	4-Bromophenyl phenyl ether	<1.2	ug/L		1.2	4	7/30/2001	JF
SW-846-8270C	4-Chloro-3-methylphenol	<1.7	ug/L		1.7	5.6666	7/30/2001	JF
SW-846-8270C	4-Chloroaniline	<1.3	ug/L		1.32	4.4	7/30/2001	JF
SW-846-8270C	4-Chlorophenyl phenyl ether	<1.2	ug/L		1.24	4.1333	7/30/2001	JF
SW-846-8270C	4-Nitroaniline	<1.3	ug/L		1.32	4.4	7/30/2001	JF
SW-846-8270C	4-Nitrophenol	<1.9	ug/L		1.86	6.2	7/30/2001	JF
SW-846-8270C	7,12-Dimethylbenz(a)anthracene	<2.0	ug/L		2	6.6666	7/30/2001	JF
SW-846-8270C	Acenaphthene	<1.1	ug/L		1.12	3.7333	7/30/2001	JF
SW-846-8270C	Acenaphthylene	<1.5	ug/L		1.46	4.8666	7/30/2001	JF
SW-846-8270C	Acetophenone	<1.6	ug/L		1.6	5.3333	7/30/2001	JF
SW-846-8270C	Aniline	<1.7	ug/L		1.74	5.8	7/30/2001	JF
SW-846-8270C	Anthracene	<1.3	ug/L		1.3	4.3333	7/30/2001	JF
SW-846-8270C	Benzidine	<2.8	ug/L		2.8	9.3333	7/30/2001	JF

Robert E. Lee & Associates, Inc.
Wisconsin Certification Number: 405043870
Certificate of Analysis Report

Robert E Lee & Associates, Inc
2825 S Webster Ave

Green Bay, WI 54301-2878

Project Number: 13551002

Project Name: WEISENBERGER TIE & LUMBER

Attn: Jim Caine

Phone: (920)336-6338

Fax: (920)336-9141

Client ID: L14

Chain: 85968

Report Date: 8/23/2001

Method	Parameter Name	Result	Units	Flag	MDL	PQL	Anl Date	Analyst
Lab No.	Collect Date	Sample ID						
SW-846-8270C		Benzo(a)anthracene	<1.4	ug/L	1.42	4.7333	7/30/2001	JF
SW-846-8270C		Benzo(a)pyrene	<1.1	ug/L	1.1	3.6666	7/30/2001	JF
SW-846-8270C		Benzo(b)fluoranthene	<1.3	ug/L	1.28	4.2666	7/30/2001	JF
SW-846-8270C		Benzo(g,h,i)perylene	<1.3	ug/L	1.32	4.4	7/30/2001	JF
SW-846-8270C		Benzo(k)fluoranthene	<1.6	ug/L	1.58	5.2666	7/30/2001	JF
SW-846-8270C		Benzyl alcohol	<1.8	ug/L	1.8	6	7/30/2001	JF
SW-846-8270C		bis(2-Chloroethoxy)methane	<1.5	ug/L	1.48	4.9333	7/30/2001	JF
SW-846-8270C		bis(2-Chloroethyl)ether	<1.9	ug/L	1.9	6.3333	7/30/2001	JF
SW-846-8270C		bis(2-Chloroisopropyl)ether	<1.5	ug/L	1.54	5.1333	7/30/2001	JF
SW-846-8270C		bis(2-Ethylhexyl)phthalate	<4.6	ug/L	4.6	15.333	7/30/2001	JF
SW-846-8270C		Butylbenzylphthalate	<1.8	ug/L	1.8	6	7/30/2001	JF
SW-846-8270C		Chrysene	<1.6	ug/L	1.56	5.2	7/30/2001	JF
SW-846-8270C		Di-n-butylphthalate	<1.5	ug/L	1.52	5.0666	7/30/2001	JF
SW-846-8270C		Di-n-octylphthalate	<1.5	ug/L	1.52	5.0666	7/30/2001	JF
SW-846-8270C		Dibenz(a,j)acridine	<1.4	ug/L	1.44	4.8	7/30/2001	JF
SW-846-8270C		Dibenzo(a,h)anthracene	<1.5	ug/L	1.5	5	7/30/2001	JF
SW-846-8270C		Dibenzofuran	<1.3	ug/L	1.28	4.2666	7/30/2001	JF
SW-846-8270C		Diethylphthalate	<1.3	ug/L	1.26	4.2	7/30/2001	JF
SW-846-8270C		Dimethylphthalate	<1.5	ug/L	1.52	5.0666	7/30/2001	JF
SW-846-8270C		Diphenylamine	<2.6	ug/L	2.6	8.6666	7/30/2001	JF
SW-846-8270C		Ethyl methanesulfonate	<1.7	ug/L	1.7	5.6666	7/30/2001	JF
SW-846-8270C		Fluoranthene	<2.0	ug/L	1.98	6.6	7/30/2001	JF
SW-846-8270C		Fluorene	<1.2	ug/L	1.2	4	7/30/2001	JF
SW-846-8270C		Hexachlorobenzene	<1.6	ug/L	1.58	5.2666	7/30/2001	JF
SW-846-8270C		Hexachlorobutadiene	<7.6	ug/L	7.6	25.333	7/30/2001	JF
SW-846-8270C		Hexachlorocyclopentadiene	<5.4	ug/L	5.4	18	7/30/2001	JF
SW-846-8270C		Hexachloroethane	<5.2	ug/L	5.2	17.333	7/30/2001	JF
SW-846-8270C		Indeno(1,2,3-cd)pyrene	<1.4	ug/L	1.44	4.8	7/30/2001	JF
SW-846-8270C		Isophorone	<1.4	ug/L	1.42	4.7333	7/30/2001	JF
SW-846-8270C		Methyl methanesulfonate	<1.1	ug/L	1.06	3.5333	7/30/2001	JF
SW-846-8270C		n-Nitrosodi-n-butylamine	<1.5	ug/L	1.46	4.8666	7/30/2001	JF
SW-846-8270C		n-Nitrosodi-n-propylamine	<1.4	ug/L	1.4	4.6666	7/30/2001	JF
SW-846-8270C		n-Nitrosodimethylamine	<1.9	ug/L	1.88	6.2666	7/30/2001	JF
SW-846-8270C		n-Nitrosodiphenylamine	<2.6	ug/L	2.6	8.6666	7/30/2001	JF
SW-846-8270C		n-Nitrosopiperidine	<1.4	ug/L	1.4	4.6666	7/30/2001	JF
SW-846-8270C		Naphthalene	<1.4	ug/L	1.38	4.6	7/30/2001	JF
SW-846-8270C		Nitrobenzene	<1.4	ug/L	1.42	4.7333	7/30/2001	JF
SW-846-8270C		Nitrobenzene-d5 - Surrogate	53	% Rec	0	0	7/30/2001	JF
SW-846-8270C		p-Dimethylaminoazobenzene	<2.0	ug/L	2	6.6666	7/30/2001	JF
SW-846-8270C		Pentachlorobenzene	<2.2	ug/L	2.2	7.3333	7/30/2001	JF
SW-846-8270C		Pentachloronitrobenzene	<1.5	ug/L	1.54	5.1333	7/30/2001	JF
SW-846-8270C		Pentachlorophenol	<1.7	ug/L	1.72	5.7333	7/30/2001	JF
SW-846-8270C		Phenacetin	<1.8	ug/L	1.78	5.9333	7/30/2001	JF
SW-846-8270C		Phenanthrene	<1.4	ug/L	1.42	4.7333	7/30/2001	JF
SW-846-8270C		Phenol	<1.1	ug/L	1.1	3.6666	7/30/2001	JF
SW-846-8270C		Phenol-d5 - Surrogate	38	% Rec	0	0	7/30/2001	JF

Robert E. Lee & Associates, Inc.
 Wisconsin Certification Number: 405043870
 Certificate of Analysis Report

Robert E Lee & Associates, Inc
 2825 S Webster Ave

Attn: Jim Caine
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 Fax: (920)336-9141

Green Bay, WI 54301-2878
 Project Number: 13551002

Client ID: L14
 Chain: 85968

Project Name: WEISENBERGER TIE & LUMBER

Report Date: 8/23/2001

Method	Parameter Name	Result	Units	Flag	MDL	PQL	Anl Date	Analyst
Lab No	Collect Date	Sample ID						
SW-846-8270C	Pronamide	<1.7	ug/L		1.68	5.6	7/30/2001	JF
SW-846-8270C	Pyrene	<2.2	ug/L		2.2	7.3333	7/30/2001	JF
SW-846-8270C	Pyridine	<3.0	ug/L		3	10	7/30/2001	JF
SW-846-8270C	Terphenyl-d14 - Surrogate	57	% Rec		0	0	7/30/2001	JF
01REL010552	6/27/2001	DMW-6A						
EPA 1613	Dioxin Analysis	See Attached			0	0	7/25/2001	PJK
SW-846-8021B	1,2,4-Trimethylbenzene	<0.23	ug/L		0.23	0.7666	7/10/2001	CRW
SW-846-8021B	1,3,5-Trimethylbenzene	<0.21	ug/L		0.21	0.7	7/10/2001	CRW
SW-846-8021B	Benzene	<0.21	ug/L		0.21	0.7	7/10/2001	CRW
SW-846-8021B	Ethylbenzene	<0.23	ug/L		0.23	0.7666	7/10/2001	CRW
SW-846-8021B	Fluorobenzene-Surrogate	100	% Rec		0	0	7/10/2001	CRW
SW-846-8021B	Methyl-tertiary-butyl ether	<0.091	ug/L		0.091	0.3033	7/10/2001	CRW
SW-846-8021B	Toluene	<0.22	ug/L		0.22	0.7333	7/10/2001	CRW
SW-846-8021B	Xylenes-Total	<0.44	ug/L		0.44	1.4666	7/10/2001	CRW
SW-846-8270C	Extraction Date	Complete			0	0	7/03/2001	JF
SW-846-8270C	1,2,4,5-Tetrachlorobenzene	<3.0	ug/L		3	10	7/30/2001	JF
SW-846-8270C	1,2,4-Trichlorobenzene	<3.6	ug/L		3.6	12	7/30/2001	JF
SW-846-8270C	1,2-Dichlorobenzene	<3.2	ug/L		3.2	10.666	7/30/2001	JF
SW-846-8270C	1,2-Diphenylhydrazine	<1.3	ug/L		1.34	4.4666	7/30/2001	JF
SW-846-8270C	1,3-Dichlorobenzene	<3.6	ug/L		3.6	12	7/30/2001	JF
SW-846-8270C	1,4-Dichlorobenzene	<3.4	ug/L		3.4	11.333	7/30/2001	JF
SW-846-8270C	1-Chloronaphthalene	<1.8	ug/L		1.84	6.1333	7/30/2001	JF
SW-846-8270C	1-Naphthylamine	<2.2	ug/L		2.2	7.3333	7/30/2001	JF
SW-846-8270C	2,3,4,6-Tetrachlorophenol	<1.4	ug/L		1.36	4.5333	7/30/2001	JF
SW-846-8270C	2,4,5-Trichlorophenol	<1.6	ug/L		1.64	5.4666	7/30/2001	JF
SW-846-8270C	2,4,6-Tribromophenol - Surrogate	63	% Rec		0	0	7/30/2001	JF
SW-846-8270C	2,4,6-Trichlorophenol	<1.6	ug/L		1.62	5.4	7/30/2001	JF
SW-846-8270C	2,4-Dichlorophenol	<1.5	ug/L		1.54	5.1333	7/30/2001	JF
SW-846-8270C	2,4-Dimethylphenol	<2.8	ug/L		2.8	9.3333	7/30/2001	JF
SW-846-8270C	2,4-Dinitrophenol	<1.8	ug/L		1.78	5.9333	7/30/2001	JF
SW-846-8270C	2,4-Dinitrotoluene	<1.6	ug/L		1.58	5.2666	7/30/2001	JF
SW-846-8270C	2,6-Dichlorophenol	<1.5	ug/L		1.52	5.0666	7/30/2001	JF
SW-846-8270C	2,6-Dinitrotoluene	<1.6	ug/L		1.6	5.3333	7/30/2001	JF
SW-846-8270C	2-Chloronaphthalene	<1.6	ug/L		1.56	5.2	7/30/2001	JF
SW-846-8270C	2-Chlorophenol	<1.2	ug/L		1.2	4	7/30/2001	JF
SW-846-8270C	2-Fluorobiphenyl - Surrogate	60	% Rec		0	0	7/30/2001	JF
SW-846-8270C	2-Fluorophenol - Surrogate	42	% Rec		0	0	7/30/2001	JF
SW-846-8270C	2-Methylnaphthalene	<1.6	ug/L		1.6	5.3333	7/30/2001	JF
SW-846-8270C	2-Methylphenol	<1.7	ug/L		1.7	5.6666	7/30/2001	JF
SW-846-8270C	2-Naphthylamine	<2.2	ug/L		2.2	7.3333	7/30/2001	JF
SW-846-8270C	2-Nitroaniline	<1.6	ug/L		1.64	5.4666	7/30/2001	JF
SW-846-8270C	2-Nitrophenol	<1.3	ug/L		1.28	4.2666	7/30/2001	JF
SW-846-8270C	2-Picoline	<3.0	ug/L		3	10	7/30/2001	JF
SW-846-8270C	3 & 4-Methylphenol	<1.8	ug/L		1.76	5.8666	7/30/2001	JF
SW-846-8270C	3,3'-Dichlorobenzidine	<0.84	ug/L		0.84	2.8	7/30/2001	JF
SW-846-8270C	3-Methylcholanthrene	<1.9	ug/L		1.94	6.4666	7/30/2001	JF

Robert E Lee & Associates, Inc
 2825 S Webster Ave

Green Bay, WI 54301-2878
 Project Number: 13551002

Project Name: WEISENBERGER TIE & LUMBER

Attn: Jim Caine

Phone: (920)336-6338

Fax: (920)336-9141

Client ID: L14

Chain: 85968

Report Date: 8/23/2001

Method	Parameter Name	Result	Units	Flag	MDL	PQL	Anl Date	Analyst
Lab No.	Collect Date	Sample ID						
SW-846-8270C	3-Nitroaniline	<2.2	ug/L		2.2	7.3333	7/30/2001	JF
SW-846-8270C	4,6-Dinitro-2-methylphenol	<2.6	ug/L		2.6	8.6666	7/30/2001	JF
SW-846-8270C	4-Aminobiphenyl	<1.5	ug/L		1.5	5	7/30/2001	JF
SW-846-8270C	4-Bromophenyl phenyl ether	<1.2	ug/L		1.2	4	7/30/2001	JF
SW-846-8270C	4-Chloro-3-methylphenol	<1.7	ug/L		1.7	5.6666	7/30/2001	JF
SW-846-8270C	4-Chloroaniline	<1.3	ug/L		1.32	4.4	7/30/2001	JF
SW-846-8270C	4-Chlorophenyl phenyl ether	<1.2	ug/L		1.24	4.1333	7/30/2001	JF
SW-846-8270C	4-Nitroaniline	<1.3	ug/L		1.32	4.4	7/30/2001	JF
SW-846-8270C	4-Nitrophenol	<1.9	ug/L		1.86	6.2	7/30/2001	JF
SW-846-8270C	7,12-Dimethylbenz(a)anthracene	<2.0	ug/L		2	6.6666	7/30/2001	JF
SW-846-8270C	Acenaphthene	<1.1	ug/L		1.12	3.7333	7/30/2001	JF
SW-846-8270C	Acenaphthylene	<1.5	ug/L		1.46	4.8666	7/30/2001	JF
SW-846-8270C	Acetophenone	<1.6	ug/L		1.6	5.3333	7/30/2001	JF
SW-846-8270C	Aniline	<1.7	ug/L		1.74	5.8	7/30/2001	JF
SW-846-8270C	Anthracene	<1.3	ug/L		1.3	4.3333	7/30/2001	JF
SW-846-8270C	Benzidine	<2.8	ug/L		2.8	9.3333	7/30/2001	JF
SW-846-8270C	Benzo(a)anthracene	<1.4	ug/L		1.42	4.7333	7/30/2001	JF
SW-846-8270C	Benzo(a)pyrene	<1.1	ug/L		1.1	3.6666	7/30/2001	JF
SW-846-8270C	Benzo(b)fluoranthene	<1.3	ug/L		1.28	4.2666	7/30/2001	JF
SW-846-8270C	Benzo(g,h,i)perylene	<1.3	ug/L		1.32	4.4	7/30/2001	JF
SW-846-8270C	Benzo(k)fluoranthene	<1.6	ug/L		1.58	5.2666	7/30/2001	JF
SW-846-8270C	Benzyl alcohol	<1.8	ug/L		1.8	6	7/30/2001	JF
SW-846-8270C	bis(2-Chloroethoxy)methane	<1.5	ug/L		1.48	4.9333	7/30/2001	JF
SW-846-8270C	bis(2-Chloroethyl)ether	<1.9	ug/L		1.9	6.3333	7/30/2001	JF
SW-846-8270C	bis(2-Chloroisopropyl)ether	<1.5	ug/L		1.54	5.1333	7/30/2001	JF
SW-846-8270C	bis(2-Ethylhexyl)phthalate	<4.6	ug/L		4.6	15.333	7/30/2001	JF
SW-846-8270C	Butylbenzylphthalate	<1.8	ug/L		1.8	6	7/30/2001	JF
SW-846-8270C	Chrysene	<1.6	ug/L		1.56	5.2	7/30/2001	JF
SW-846-8270C	Di-n-butylphthalate	<1.5	ug/L		1.52	5.0666	7/30/2001	JF
SW-846-8270C	Di-n-octylphthalate	<1.5	ug/L		1.52	5.0666	7/30/2001	JF
SW-846-8270C	Dibenz(a,j)acridine	<1.4	ug/L		1.44	4.8	7/30/2001	JF
SW-846-8270C	Dibenzo(a,h)anthracene	<1.5	ug/L		1.5	5	7/30/2001	JF
SW-846-8270C	Dibenzofuran	<1.3	ug/L		1.28	4.2666	7/30/2001	JF
SW-846-8270C	Diethylphthalate	<1.3	ug/L		1.26	4.2	7/30/2001	JF
SW-846-8270C	Dimethylphthalate	<1.5	ug/L		1.52	5.0666	7/30/2001	JF
SW-846-8270C	Diphenylamine	<2.6	ug/L		2.6	8.6666	7/30/2001	JF
SW-846-8270C	Ethyl methanesulfonate	<1.7	ug/L		1.7	5.6666	7/30/2001	JF
SW-846-8270C	Fluoranthene	<2.0	ug/L		1.98	6.6	7/30/2001	JF
SW-846-8270C	Fluorene	<1.2	ug/L		1.2	4	7/30/2001	JF
SW-846-8270C	Hexachlorobenzene	<1.6	ug/L		1.58	5.2666	7/30/2001	JF
SW-846-8270C	Hexachlorobutadiene	<7.6	ug/L		7.6	25.333	7/30/2001	JF
SW-846-8270C	Hexachlorocyclopentadiene	<5.4	ug/L		5.4	18	7/30/2001	JF
SW-846-8270C	Hexachloroethane	<5.2	ug/L		5.2	17.333	7/30/2001	JF
SW-846-8270C	Indeno(1,2,3-cd)pyrene	<1.4	ug/L		1.44	4.8	7/30/2001	JF
SW-846-8270C	Isophorone	<1.4	ug/L		1.42	4.7333	7/30/2001	JF
SW-846-8270C	Methyl methanesulfonate	<1.1	ug/L		1.06	3.5333	7/30/2001	JF

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Method	Parameter Name	Result	Units	Flag	MDL	PQL	Anl Date	Analyst
Lab No	Collect Date	Sample ID						
SW-846-8270C	n-Nitrosodi-n-butylamine	<1.5	ug/L		1.46	4.8666	7/30/2001	JF
SW-846-8270C	n-Nitrosodi-n-propylamine	<1.4	ug/L		1.4	4.6666	7/30/2001	JF
SW-846-8270C	n-Nitrosodimethylamine	<1.9	ug/L		1.88	6.2666	7/30/2001	JF
SW-846-8270C	n-Nitrosodiphenylamine	<2.6	ug/L		2.6	8.6666	7/30/2001	JF
SW-846-8270C	n-Nitrosopiperidine	<1.4	ug/L		1.4	4.6666	7/30/2001	JF
SW-846-8270C	Naphthalene	<1.4	ug/L		1.38	4.6	7/30/2001	JF
SW-846-8270C	Nitrobenzene	<1.4	ug/L		1.42	4.7333	7/30/2001	JF
SW-846-8270C	Nitrobenzene-d5 - Surrogate	52	% Rec		0	0	7/30/2001	JF
SW-846-8270C	p-Dimethylaminoazobenzene	<2.0	ug/L		2	6.6666	7/30/2001	JF
SW-846-8270C	Pentachlorobenzene	<2.2	ug/L		2.2	7.3333	7/30/2001	JF
SW-846-8270C	Pentachloronitrobenzene	<1.5	ug/L		1.54	5.1333	7/30/2001	JF
SW-846-8270C	Pentachlorophenol	<1.7	ug/L		1.72	5.7333	7/30/2001	JF
SW-846-8270C	Phenacetin	<1.8	ug/L		1.78	5.9333	7/30/2001	JF
SW-846-8270C	Phenanthrene	<1.4	ug/L		1.42	4.7333	7/30/2001	JF
SW-846-8270C	Phenol	<1.1	ug/L		1.1	3.6666	7/30/2001	JF
SW-846-8270C	Phenol-d5 - Surrogate	37	% Rec		0	0	7/30/2001	JF
SW-846-8270C	Pronamide	<1.7	ug/L		1.68	5.6	7/30/2001	JF
SW-846-8270C	Pyrene	<2.2	ug/L		2.2	7.3333	7/30/2001	JF
SW-846-8270C	Pyridine	<3.0	ug/L		3	10	7/30/2001	JF
SW-846-8270C	Terphenyl-d14 - Surrogate	58	% Rec		0	0	7/30/2001	JF

01RELO10553 6/27/2001 DMW-7

EPA 1613	Dioxin Analysis	See Attached			0	0	7/25/2001	PJK
SW-846-8270C	Extraction Date	Complete			0	0	7/03/2001	JF
SW-846-8270C	1,2,4,5-Tetrachlorobenzene	<3.0	ug/L		3	10	7/30/2001	JF
SW-846-8270C	1,2,4-Trichlorobenzene	<3.6	ug/L		3.6	12	7/30/2001	JF
SW-846-8270C	1,2-Dichlorobenzene	<3.2	ug/L		3.2	10.666	7/30/2001	JF
SW-846-8270C	1,2-Diphenylhydrazine	<1.3	ug/L		1.34	4.4666	7/30/2001	JF
SW-846-8270C	1,3-Dichlorobenzene	<3.6	ug/L		3.6	12	7/30/2001	JF
SW-846-8270C	1,4-Dichlorobenzene	<3.4	ug/L		3.4	11.333	7/30/2001	JF
SW-846-8270C	1-Chloronaphthalene	<1.8	ug/L		1.84	6.1333	7/30/2001	JF
SW-846-8270C	1-Naphthylamine	<2.2	ug/L		2.2	7.3333	7/30/2001	JF
SW-846-8270C	2,3,4,6-Tetrachlorophenol	<1.4	ug/L		1.36	4.5333	7/30/2001	JF
SW-846-8270C	2,4,5-Trichlorophenol	<1.6	ug/L		1.64	5.4666	7/30/2001	JF
SW-846-8270C	2,4,6-Tribromophenol - Surrogate	64	% Rec		0	0	7/30/2001	JF
SW-846-8270C	2,4,6-Trichlorophenol	<1.6	ug/L		1.62	5.4	7/30/2001	JF
SW-846-8270C	2,4-Dichlorophenol	<1.5	ug/L		1.54	5.1333	7/30/2001	JF
SW-846-8270C	2,4-Dimethylphenol	<2.8	ug/L		2.8	9.3333	7/30/2001	JF
SW-846-8270C	2,4-Dinitrophenol	<1.8	ug/L		1.78	5.9333	7/30/2001	JF
SW-846-8270C	2,4-Dinitrotoluene	<1.6	ug/L		1.58	5.2666	7/30/2001	JF
SW-846-8270C	2,6-Dichlorophenol	<1.5	ug/L		1.52	5.0666	7/30/2001	JF
SW-846-8270C	2,6-Dinitrotoluene	<1.6	ug/L		1.6	5.3333	7/30/2001	JF
SW-846-8270C	2-Chloronaphthalene	<1.6	ug/L		1.56	5.2	7/30/2001	JF
SW-846-8270C	2-Chlorophenol	<1.2	ug/L		1.2	4	7/30/2001	JF
SW-846-8270C	2-Fluorobiphenyl - Surrogate	62	% Rec		0	0	7/30/2001	JF
SW-846-8270C	2-Fluorophenol - Surrogate	44	% Rec		0	0	7/30/2001	JF
SW-846-8270C	2-Methylnaphthalene	<1.6	ug/L		1.6	5.3333	7/30/2001	JF

Robert E Lee & Associates, Inc
 2825 S Webster Ave

Green Bay, WI 54301-2878
 Project Number: 13551002
 Project Name: WEISENBERGER TIE & LUMBER

Attn: Jim Caine
 Phone: (920)336-6338
 Fax: (920)336-9141
 Client ID: L14
 Chain: 85968
 Report Date: 8/23/2001

Method	Parameter Name	Result	Units	Flag	MDL	PQL	Anl Date	Analyst
Lab No	Collect Date	Sample ID						
SW-846-8270C	2-Methylphenol	<1.7	ug/L		1.7	5.6666	7/30/2001	JF
SW-846-8270C	2-Naphthylamine	<2.2	ug/L		2.2	7.3333	7/30/2001	JF
SW-846-8270C	2-Nitroaniline	<1.6	ug/L		1.64	5.4666	7/30/2001	JF
SW-846-8270C	2-Nitrophenol	<1.3	ug/L		1.28	4.2666	7/30/2001	JF
SW-846-8270C	2-Picoline	<3.0	ug/L		3	10	7/30/2001	JF
SW-846-8270C	3 & 4-Methylphenol	<1.8	ug/L		1.76	5.8666	7/30/2001	JF
SW-846-8270C	3,3'-Dichlorobenzidine	<0.84	ug/L		0.84	2.8	7/30/2001	JF
SW-846-8270C	3-Methylcholanthrene	<1.9	ug/L		1.94	6.4666	7/30/2001	JF
SW-846-8270C	3-Nitroaniline	<2.2	ug/L		2.2	7.3333	7/30/2001	JF
SW-846-8270C	4,6-Dinitro-2-methylphenol	<2.6	ug/L		2.6	8.6666	7/30/2001	JF
SW-846-8270C	4-Aminobiphenyl	<1.5	ug/L		1.5	5	7/30/2001	JF
SW-846-8270C	4-Bromophenyl phenyl ether	<1.2	ug/L		1.2	4	7/30/2001	JF
SW-846-8270C	4-Chloro-3-methylphenol	<1.7	ug/L		1.7	5.6666	7/30/2001	JF
SW-846-8270C	4-Chloroaniline	<1.3	ug/L		1.32	4.4	7/30/2001	JF
SW-846-8270C	4-Chlorophenyl phenyl ether	<1.2	ug/L		1.24	4.1333	7/30/2001	JF
SW-846-8270C	4-Nitroaniline	<1.3	ug/L		1.32	4.4	7/30/2001	JF
SW-846-8270C	4-Nitrophenol	<1.9	ug/L		1.86	6.2	7/30/2001	JF
SW-846-8270C	7,12-Dimethylbenz(a)anthracene	<2.0	ug/L		2	6.6666	7/30/2001	JF
SW-846-8270C	Acenaphthene	<1.1	ug/L		1.12	3.7333	7/30/2001	JF
SW-846-8270C	Acenaphthylene	<1.5	ug/L		1.46	4.8666	7/30/2001	JF
SW-846-8270C	Acetophenone	<1.6	ug/L		1.6	5.3333	7/30/2001	JF
SW-846-8270C	Aniline	<1.7	ug/L		1.74	5.8	7/30/2001	JF
SW-846-8270C	Anthracene	<1.3	ug/L		1.3	4.3333	7/30/2001	JF
SW-846-8270C	Benzidine	<2.8	ug/L		2.8	9.3333	7/30/2001	JF
SW-846-8270C	Benzo(a)anthracene	<1.4	ug/L		1.42	4.7333	7/30/2001	JF
SW-846-8270C	Benzo(a)pyrene	<1.1	ug/L		1.1	3.6666	7/30/2001	JF
SW-846-8270C	Benzo(b)fluoranthene	<1.3	ug/L		1.28	4.2666	7/30/2001	JF
SW-846-8270C	Benzo(g,h,i)perylene	<1.3	ug/L		1.32	4.4	7/30/2001	JF
SW-846-8270C	Benzo(k)fluoranthene	<1.6	ug/L		1.58	5.2666	7/30/2001	JF
SW-846-8270C	Benzyl alcohol	<1.8	ug/L		1.8	6	7/30/2001	JF
SW-846-8270C	bis(2-Chloroethoxy)methane	<1.5	ug/L		1.48	4.9333	7/30/2001	JF
SW-846-8270C	bis(2-Chloroethyl)ether	<1.9	ug/L		1.9	6.3333	7/30/2001	JF
SW-846-8270C	bis(2-Chloroisopropyl)ether	<1.5	ug/L		1.54	5.1333	7/30/2001	JF
SW-846-8270C	bis(2-Ethylhexyl)phthalate	<4.6	ug/L		4.6	15.333	7/30/2001	JF
SW-846-8270C	Butylbenzylphthalate	<1.8	ug/L		1.8	6	7/30/2001	JF
SW-846-8270C	Chrysene	<1.6	ug/L		1.56	5.2	7/30/2001	JF
SW-846-8270C	Di-n-butylphthalate	<1.5	ug/L		1.52	5.0666	7/30/2001	JF
SW-846-8270C	Di-n-octylphthalate	<1.5	ug/L		1.52	5.0666	7/30/2001	JF
SW-846-8270C	Dibenz(a,j)acridine	<1.4	ug/L		1.44	4.8	7/30/2001	JF
SW-846-8270C	Dibenzo(a,h)anthracene	<1.5	ug/L		1.5	5	7/30/2001	JF
SW-846-8270C	Dibenzofuran	<1.3	ug/L		1.28	4.2666	7/30/2001	JF
SW-846-8270C	Diethylphthalate	<1.3	ug/L		1.26	4.2	7/30/2001	JF
SW-846-8270C	Dimethylphthalate	<1.5	ug/L		1.52	5.0666	7/30/2001	JF
SW-846-8270C	Diphenylamine	<2.6	ug/L		2.6	8.6666	7/30/2001	JF
SW-846-8270C	Ethyl methanesulfonate	<1.7	ug/L		1.7	5.6666	7/30/2001	JF
SW-846-8270C	Fluoranthene	<2.0	ug/L		1.98	6.6	7/30/2001	JF

Robert E. Lee & Associates, Inc.
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Certificate of Analysis Report

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Green Bay, WI 54301-2878
Project Number: 13551002
Project Name: WEISENBERGER TIE & LUMBER

Attn: Jim Caine
Phone: (920)336-6338
Fax: (920)336-9141
Client ID: L14
Chain: 85968
Report Date: 8/23/2001

Method	Parameter Name	Result	Units	Flag	MDL	PQL	Anl Date	Analyst
Lab No	Collect Date	Sample ID						
SW-846-8270C	Fluorene	<1.2	ug/L		1.2	4	7/30/2001	JF
SW-846-8270C	Hexachlorobenzene	<1.6	ug/L		1.58	5.2666	7/30/2001	JF
SW-846-8270C	Hexachlorobutadiene	<7.6	ug/L		7.6	25.333	7/30/2001	JF
SW-846-8270C	Hexachlorocyclopentadiene	<5.4	ug/L		5.4	18	7/30/2001	JF
SW-846-8270C	Hexachloroethane	<5.2	ug/L		5.2	17.333	7/30/2001	JF
SW-846-8270C	Indeno(1,2,3-cd)pyrene	<1.4	ug/L		1.44	4.8	7/30/2001	JF
SW-846-8270C	Isophorone	<1.4	ug/L		1.42	4.7333	7/30/2001	JF
SW-846-8270C	Methyl methanesulfonate	<1.1	ug/L		1.06	3.5333	7/30/2001	JF
SW-846-8270C	n-Nitrosodi-n-butylamine	<1.5	ug/L		1.46	4.8666	7/30/2001	JF
SW-846-8270C	n-Nitrosodi-n-propylamine	<1.4	ug/L		1.4	4.6666	7/30/2001	JF
SW-846-8270C	n-Nitrosodimethylamine	<1.9	ug/L		1.88	6.2666	7/30/2001	JF
SW-846-8270C	n-Nitrosodiphenylamine	<2.6	ug/L		2.6	8.6666	7/30/2001	JF
SW-846-8270C	n-Nitrosopiperidine	<1.4	ug/L		1.4	4.6666	7/30/2001	JF
SW-846-8270C	Naphthalene	<1.4	ug/L		1.38	4.6	7/30/2001	JF
SW-846-8270C	Nitrobenzene	<1.4	ug/L		1.42	4.7333	7/30/2001	JF
SW-846-8270C	Nitrobenzene-d5 - Surrogate	54	% Rec		0	0	7/30/2001	JF
SW-846-8270C	p-Dimethylaminoazobenzene	<2.0	ug/L		2	6.6666	7/30/2001	JF
SW-846-8270C	Pentachlorobenzene	<2.2	ug/L		2.2	7.3333	7/30/2001	JF
SW-846-8270C	Pentachloronitrobenzene	<1.5	ug/L		1.54	5.1333	7/30/2001	JF
SW-846-8270C	Pentachlorophenol	<1.7	ug/L		1.72	5.7333	7/30/2001	JF
SW-846-8270C	Phenacetin	<1.8	ug/L		1.78	5.9333	7/30/2001	JF
SW-846-8270C	Phenanthrene	<1.4	ug/L		1.42	4.7333	7/30/2001	JF
SW-846-8270C	Phenol	<1.1	ug/L		1.1	3.6666	7/30/2001	JF
SW-846-8270C	Phenol-d5 - Surrogate	38	% Rec		0	0	7/30/2001	JF
SW-846-8270C	Pronamide	<1.7	ug/L		1.68	5.6	7/30/2001	JF
SW-846-8270C	Pyrene	<2.2	ug/L		2.2	7.3333	7/30/2001	JF
SW-846-8270C	Pyridine	<3.0	ug/L		3	10	7/30/2001	JF
SW-846-8270C	Terphenyl-d14 - Surrogate	59	% Rec		0	0	7/30/2001	JF
01REL010554	6/27/2001	DMW-8						
EPA 1613	Dioxin Analysis	See Attached			0	0	7/25/2001	PJK
01REL010555	6/27/2001	DMW-10						
SW-846-8021B	1,2,4-Trimethylbenzene	<0.23	ug/L		0.23	0.7666	7/10/2001	CRW
SW-846-8021B	1,3,5-Trimethylbenzene	<0.21	ug/L		0.21	0.7	7/10/2001	CRW
SW-846-8021B	Benzene	<0.21	ug/L		0.21	0.7	7/10/2001	CRW
SW-846-8021B	Ethylbenzene	<0.23	ug/L		0.23	0.7666	7/10/2001	CRW
SW-846-8021B	Fluorobenzene-Surrogate	103	% Rec		0	0	7/10/2001	CRW
SW-846-8021B	Methyl-tertiary-butyl ether	<0.091	ug/L		0.091	0.3033	7/10/2001	CRW
SW-846-8021B	Toluene	<0.22	ug/L		0.22	0.7333	7/10/2001	CRW
SW-846-8021B	Xylenes-Total	<0.44	ug/L		0.44	1.4666	7/10/2001	CRW
SW-846-8270C	Extraction Date	Complete			0	0	7/03/2001	JF
SW-846-8270C	1,2,4,5-Tetrachlorobenzene	<3.0	ug/L		3	10	8/09/2001	JF
SW-846-8270C	1,2,4-Trichlorobenzene	<3.6	ug/L		3.6	12	8/09/2001	JF
SW-846-8270C	1,2-Dichlorobenzene	<3.2	ug/L		3.2	10.666	8/09/2001	JF
SW-846-8270C	1,2-Diphenylhydrazine	<1.3	ug/L		1.34	4.4666	8/09/2001	JF
SW-846-8270C	1,3-Dichlorobenzene	<3.6	ug/L		3.6	12	8/09/2001	JF

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Project Name: WEISENBERGER TIE & LUMBER

Attn: Jim Caine

Phone: (920)336-6338

Fax: (920)336-9141

Client ID: L14

Chain: 85968

Report Date: 8/23/2001

Method	Parameter Name	Result	Units	Flag	MDL	PQL	Anl Date	Analyst
Lab No	Collect Date	Sample ID						
SW-846-8270C	1,4-Dichlorobenzene	<3.4	ug/L		3.4	11.333	8/09/2001	JF
SW-846-8270C	1-Chloronaphthalene	<1.8	ug/L		1.84	6.1333	8/09/2001	JF
SW-846-8270C	1-Naphthylamine	<2.2	ug/L		2.2	7.3333	8/09/2001	JF
SW-846-8270C	2,3,4,6-Tetrachlorophenol	<1.4	ug/L		1.36	4.5333	8/09/2001	JF
SW-846-8270C	2,4,5-Trichlorophenol	<1.6	ug/L		1.64	5.4666	8/09/2001	JF
SW-846-8270C	2,4,6-Tribromophenol - Surrogate	75	% Rec		0	0	8/09/2001	JF
SW-846-8270C	2,4,6-Trichlorophenol	<1.6	ug/L		1.62	5.4	8/09/2001	JF
SW-846-8270C	2,4-Dichlorophenol	<1.5	ug/L		1.54	5.1333	8/09/2001	JF
SW-846-8270C	2,4-Dimethylphenol	<2.8	ug/L		2.8	9.3333	8/09/2001	JF
SW-846-8270C	2,4-Dinitrophenol	<1.8	ug/L		1.78	5.9333	8/09/2001	JF
SW-846-8270C	2,4-Dinitrotoluene	<1.6	ug/L		1.58	5.2666	8/09/2001	JF
SW-846-8270C	2,6-Dichlorophenol	<1.5	ug/L		1.52	5.0666	8/09/2001	JF
SW-846-8270C	2,6-Dinitrotoluene	<1.6	ug/L		1.6	5.3333	8/09/2001	JF
SW-846-8270C	2-Chloronaphthalene	<1.6	ug/L		1.56	5.2	8/09/2001	JF
SW-846-8270C	2-Chlorophenol	<1.2	ug/L		1.2	4	8/09/2001	JF
SW-846-8270C	2-Fluorobiphenyl - Surrogate	68	% Rec		0	0	8/09/2001	JF
SW-846-8270C	2-Fluorophenol - Surrogate	54	% Rec		0	0	8/09/2001	JF
SW-846-8270C	2-Methylnaphthalene	<1.6	ug/L		1.6	5.3333	8/09/2001	JF
SW-846-8270C	2-Methylphenol	<1.7	ug/L		1.7	5.6666	8/09/2001	JF
SW-846-8270C	2-Naphthylamine	<2.2	ug/L		2.2	7.3333	8/09/2001	JF
SW-846-8270C	2-Nitroaniline	<1.6	ug/L		1.64	5.4666	8/09/2001	JF
SW-846-8270C	2-Nitrophenol	<1.3	ug/L		1.28	4.2666	8/09/2001	JF
SW-846-8270C	2-Picoline	<3.0	ug/L		3	10	8/09/2001	JF
SW-846-8270C	3 & 4-Methylphenol	<1.8	ug/L		1.76	5.8666	8/09/2001	JF
SW-846-8270C	3,3'-Dichlorobenzidine	<0.84	ug/L		0.84	2.8	8/09/2001	JF
SW-846-8270C	3-Methylcholanthrene	<1.9	ug/L		1.94	6.4666	8/09/2001	JF
SW-846-8270C	3-Nitroaniline	<2.2	ug/L		2.2	7.3333	8/09/2001	JF
SW-846-8270C	4,6-Dinitro-2-methylphenol	<2.6	ug/L		2.6	8.6666	8/09/2001	JF
SW-846-8270C	4-Aminobiphenyl	<1.5	ug/L		1.5	5	8/09/2001	JF
SW-846-8270C	4-Bromophenyl phenyl ether	<1.2	ug/L		1.2	4	8/09/2001	JF
SW-846-8270C	4-Chloro-3-methylphenol	<1.7	ug/L		1.7	5.6666	8/09/2001	JF
SW-846-8270C	4-Chloroaniline	<1.3	ug/L		1.32	4.4	8/09/2001	JF
SW-846-8270C	4-Chlorophenyl phenyl ether	<1.2	ug/L		1.24	4.1333	8/09/2001	JF
SW-846-8270C	4-Nitroaniline	<1.3	ug/L		1.32	4.4	8/09/2001	JF
SW-846-8270C	4-Nitrophenol	<1.9	ug/L		1.86	6.2	8/09/2001	JF
SW-846-8270C	7,12-Dimethylbenz(a)anthracene	<2.0	ug/L		2	6.6666	8/09/2001	JF
SW-846-8270C	Acenaphthene	<1.1	ug/L		1.12	3.7333	8/09/2001	JF
SW-846-8270C	Acenaphthylene	<1.5	ug/L		1.46	4.8666	8/09/2001	JF
SW-846-8270C	Acetophenone	<1.6	ug/L		1.6	5.3333	8/09/2001	JF
SW-846-8270C	Aniline	<1.7	ug/L		1.74	5.8	8/09/2001	JF
SW-846-8270C	Anthracene	<1.3	ug/L		1.3	4.3333	8/09/2001	JF
SW-846-8270C	Benzidine	<2.8	ug/L		2.8	9.3333	8/09/2001	JF
SW-846-8270C	Benzo(a)anthracene	<1.4	ug/L		1.42	4.7333	8/09/2001	JF
SW-846-8270C	Benzo(a)pyrene	<1.1	ug/L		1.1	3.6666	8/09/2001	JF
SW-846-8270C	Benzo(b)fluoranthene	<1.3	ug/L		1.28	4.2666	8/09/2001	JF
SW-846-8270C	Benzo(g,h,i)perylene	<1.3	ug/L		1.32	4.4	8/09/2001	JF

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 Chain: 85968
 Report Date: 8/23/2001

Method	Parameter Name	Result	Units	Flag	MDL	PQL	Anl Date	Analyst
Lab No.	Collect Date	Sample ID						
SW-846-8270C		Benzo(k)fluoranthene	<1.6	ug/L	1.58	5.2666	8/09/2001	JF
SW-846-8270C		Benzyl alcohol	<1.8	ug/L	1.8	6	8/09/2001	JF
SW-846-8270C		bis(2-Chloroethoxy)methane	<1.5	ug/L	1.48	4.9333	8/09/2001	JF
SW-846-8270C		bis(2-Chloroethyl)ether	<1.9	ug/L	1.9	6.3333	8/09/2001	JF
SW-846-8270C		bis(2-Chloroisopropyl)ether	<1.5	ug/L	1.54	5.1333	8/09/2001	JF
SW-846-8270C		bis(2-Ethylhexyl)phthalate	<4.6	ug/L	4.6	15.333	8/09/2001	JF
SW-846-8270C		Butylbenzylphthalate	<1.8	ug/L	1.8	6	8/09/2001	JF
SW-846-8270C		Chrysene	<1.6	ug/L	1.56	5.2	8/09/2001	JF
SW-846-8270C		Di-n-butylphthalate	2.0	ug/L	13	5.0666	8/09/2001	JF
SW-846-8270C		Di-n-octylphthalate	<1.5	ug/L	1.52	5.0666	8/09/2001	JF
SW-846-8270C		Dibenz(a,j)acridine	<1.4	ug/L	1.44	4.8	8/09/2001	JF
SW-846-8270C		Dibenzo(a,h)anthracene	<1.5	ug/L	1.5	5	8/09/2001	JF
SW-846-8270C		Dibenzofuran	<1.3	ug/L	1.28	4.2666	8/09/2001	JF
SW-846-8270C		Diethylphthalate	<1.3	ug/L	1.26	4.2	8/09/2001	JF
SW-846-8270C		Dimethylphthalate	<1.5	ug/L	1.52	5.0666	8/09/2001	JF
SW-846-8270C		Diphenylamine	<2.6	ug/L	2.6	8.6666	8/09/2001	JF
SW-846-8270C		Ethyl methanesulfonate	<1.7	ug/L	1.7	5.6666	8/09/2001	JF
SW-846-8270C		Fluoranthene	<2.0	ug/L	1.98	6.6	8/09/2001	JF
SW-846-8270C		Fluorene	<1.2	ug/L	1.2	4	8/09/2001	JF
SW-846-8270C		Hexachlorobenzene	<1.6	ug/L	1.58	5.2666	8/09/2001	JF
SW-846-8270C		Hexachlorobutadiene	<7.6	ug/L	7.6	25.333	8/09/2001	JF
SW-846-8270C		Hexachlorocyclopentadiene	<5.4	ug/L	5.4	18	8/09/2001	JF
SW-846-8270C		Hexachloroethane	<5.2	ug/L	5.2	17.333	8/09/2001	JF
SW-846-8270C		Indeno(1,2,3-cd)pyrene	<1.4	ug/L	1.44	4.8	8/09/2001	JF
SW-846-8270C		Isophorone	<1.4	ug/L	1.42	4.7333	8/09/2001	JF
SW-846-8270C		Methyl methanesulfonate	<1.1	ug/L	1.06	3.5333	8/09/2001	JF
SW-846-8270C		n-Nitrosodi-n-butylamine	<1.5	ug/L	1.46	4.8666	8/09/2001	JF
SW-846-8270C		n-Nitrosodi-n-propylamine	<1.4	ug/L	1.4	4.6666	8/09/2001	JF
SW-846-8270C		n-Nitrosodimethylamine	<1.9	ug/L	1.88	6.2666	8/09/2001	JF
SW-846-8270C		n-Nitrosodiphenylamine	<2.6	ug/L	2.6	8.6666	8/09/2001	JF
SW-846-8270C		n-Nitrosopiperidine	<1.4	ug/L	1.4	4.6666	8/09/2001	JF
SW-846-8270C		Naphthalene	<1.4	ug/L	1.38	4.6	8/09/2001	JF
SW-846-8270C		Nitrobenzene	<1.4	ug/L	1.42	4.7333	8/09/2001	JF
SW-846-8270C		Nitrobenzene-d5 - Surrogate	66	% Rec	0	0	8/09/2001	JF
SW-846-8270C		p-Dimethylaminoazobenzene	<2.0	ug/L	2	6.6666	8/09/2001	JF
SW-846-8270C		Pentachlorobenzene	<2.2	ug/L	2.2	7.3333	8/09/2001	JF
SW-846-8270C		Pentachloronitrobenzene	<1.5	ug/L	1.54	5.1333	8/09/2001	JF
SW-846-8270C		Pentachlorophenol	<1.7	ug/L	1.72	5.7333	8/09/2001	JF
SW-846-8270C		Phenacetin	<1.8	ug/L	1.78	5.9333	8/09/2001	JF
SW-846-8270C		Phenanthrene	<1.4	ug/L	1.42	4.7333	8/09/2001	JF
SW-846-8270C		Phenol	<1.1	ug/L	1.1	3.6666	8/09/2001	JF
SW-846-8270C		Phenol-d5 - Surrogate	48	% Rec	0	0	8/09/2001	JF
SW-846-8270C		Pronamide	<1.7	ug/L	1.68	5.6	8/09/2001	JF
SW-846-8270C		Pyrene	<2.2	ug/L	2.2	7.3333	8/09/2001	JF
SW-846-8270C		Pyridine	<3.0	ug/L	3	10	8/09/2001	JF
SW-846-8270C		Terphenyl-d14 - Surrogate	60	% Rec	0	0	8/09/2001	JF

Robert E. Lee & Associates, Inc.
Wisconsin Certification Number: 405043870
Certificate of Analysis Report

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Project Number: 13551002
Project Name: WEISENBERGER TIE & LUMBER

Attn: Jim Caine
Phone: (920)336-6338
Fax: (920)336-9141
Client ID: L14
Chain: 85968
Report Date: 8/23/2001

Method	Parameter Name	Result	Units	Flag	MDL	PQL	Anl Date	Analyst
Lab No.	Collect Date	Sample ID						
01REL010556	6/27/2001	DMW-13						
SW-846-8270C	Extraction Date	Complete			0	0	7/03/2001	JF
SW-846-8270C	1,2,4,5-Tetrachlorobenzene	<3.0	ug/L		3	10	7/30/2001	JF
SW-846-8270C	1,2,4-Trichlorobenzene	<3.6	ug/L		3.6	12	7/30/2001	JF
SW-846-8270C	1,2-Dichlorobenzene	<3.2	ug/L		3.2	10.666	7/30/2001	JF
SW-846-8270C	1,2-Diphenylhydrazine	<1.3	ug/L		1.34	4.4666	7/30/2001	JF
SW-846-8270C	1,3-Dichlorobenzene	<3.6	ug/L		3.6	12	7/30/2001	JF
SW-846-8270C	1,4-Dichlorobenzene	<3.4	ug/L		3.4	11.333	7/30/2001	JF
SW-846-8270C	1-Chloronaphthalene	<1.8	ug/L		1.84	6.1333	7/30/2001	JF
SW-846-8270C	1-Naphthylamine	<2.2	ug/L		2.2	7.3333	7/30/2001	JF
SW-846-8270C	2,3,4,6-Tetrachlorophenol	<1.4	ug/L		1.36	4.5333	7/30/2001	JF
SW-846-8270C	2,4,5-Trichlorophenol	<1.6	ug/L		1.64	5.4666	7/30/2001	JF
SW-846-8270C	2,4,6-Tribromophenol - Surrogate	58	% Rec		0	0	7/30/2001	JF
SW-846-8270C	2,4,6-Trichlorophenol	<1.6	ug/L		1.62	5.4	7/30/2001	JF
SW-846-8270C	2,4-Dichlorophenol	<1.5	ug/L		1.54	5.1333	7/30/2001	JF
SW-846-8270C	2,4-Dimethylphenol	<2.8	ug/L		2.8	9.3333	7/30/2001	JF
SW-846-8270C	2,4-Dinitrophenol	<1.8	ug/L		1.78	5.9333	7/30/2001	JF
SW-846-8270C	2,4-Dinitrotoluene	<1.6	ug/L		1.58	5.2666	7/30/2001	JF
SW-846-8270C	2,6-Dichlorophenol	<1.5	ug/L		1.52	5.0666	7/30/2001	JF
SW-846-8270C	2,6-Dinitrotoluene	<1.6	ug/L		1.6	5.3333	7/30/2001	JF
SW-846-8270C	2-Chloronaphthalene	<1.6	ug/L		1.56	5.2	7/30/2001	JF
SW-846-8270C	2-Chlorophenol	<1.2	ug/L		1.2	4	7/30/2001	JF
SW-846-8270C	2-Fluorobiphenyl - Surrogate	55	% Rec		0	0	7/30/2001	JF
SW-846-8270C	2-Fluorophenol - Surrogate	41	% Rec		0	0	7/30/2001	JF
SW-846-8270C	2-Methylnaphthalene	<1.6	ug/L		1.6	5.3333	7/30/2001	JF
SW-846-8270C	2-Methylphenol	<1.7	ug/L		1.7	5.6666	7/30/2001	JF
SW-846-8270C	2-Naphthylamine	<2.2	ug/L		2.2	7.3333	7/30/2001	JF
SW-846-8270C	2-Nitroaniline	<1.6	ug/L		1.64	5.4666	7/30/2001	JF
SW-846-8270C	2-Nitrophenol	<1.3	ug/L		1.28	4.2666	7/30/2001	JF
SW-846-8270C	2-Picoline	<3.0	ug/L		3	10	7/30/2001	JF
SW-846-8270C	3 & 4-Methylphenol	<1.8	ug/L		1.76	5.8666	7/30/2001	JF
SW-846-8270C	3,3'-Dichlorobenzidine	<0.84	ug/L		0.84	2.8	7/30/2001	JF
SW-846-8270C	3-Methylcholanthrene	<1.9	ug/L		1.94	6.4666	7/30/2001	JF
SW-846-8270C	3-Nitroaniline	<2.2	ug/L		2.2	7.3333	7/30/2001	JF
SW-846-8270C	4,6-Dinitro-2-methylphenol	<2.6	ug/L		2.6	8.6666	7/30/2001	JF
SW-846-8270C	4-Aminobiphenyl	<1.5	ug/L		1.5	5	7/30/2001	JF
SW-846-8270C	4-Bromophenyl phenyl ether	<1.2	ug/L		1.2	4	7/30/2001	JF
SW-846-8270C	4-Chloro-3-methylphenol	<1.7	ug/L		1.7	5.6666	7/30/2001	JF
SW-846-8270C	4-Chloroaniline	<1.3	ug/L		1.32	4.4	7/30/2001	JF
SW-846-8270C	4-Chlorophenyl phenyl ether	<1.2	ug/L		1.24	4.1333	7/30/2001	JF
SW-846-8270C	4-Nitroaniline	<1.3	ug/L		1.32	4.4	7/30/2001	JF
SW-846-8270C	4-Nitrophenol	<1.9	ug/L		1.86	6.2	7/30/2001	JF
SW-846-8270C	7,12-Dimethylbenz(a)anthracene	<2.0	ug/L		2	6.6666	7/30/2001	JF
SW-846-8270C	Acenaphthene	<1.1	ug/L		1.12	3.7333	7/30/2001	JF
SW-846-8270C	Acenaphthylene	<1.5	ug/L		1.46	4.8666	7/30/2001	JF
SW-846-8270C	Acetophenone	<1.6	ug/L		1.6	5.3333	7/30/2001	JF

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 Fax: (920)336-9141
 Client ID: L14
 Chain: 85968
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Method	Parameter Name	Result	Units	Flag	MDL	PQL	Anl Date	Analyst
Lab No.	Collect Date	Sample ID						
SW-846-8270C	Aniline	<1.7	ug/L		1.74	5.8	7/30/2001	JF
SW-846-8270C	Anthracene	<1.3	ug/L		1.3	4.3333	7/30/2001	JF
SW-846-8270C	Benzidine	<2.8	ug/L		2.8	9.3333	7/30/2001	JF
SW-846-8270C	Benzo(a)anthracene	<1.4	ug/L		1.42	4.7333	7/30/2001	JF
SW-846-8270C	Benzo(a)pyrene	<1.1	ug/L		1.1	3.6666	7/30/2001	JF
SW-846-8270C	Benzo(b)fluoranthene	<1.3	ug/L		1.28	4.2666	7/30/2001	JF
SW-846-8270C	Benzo(g,h,i)perylene	<1.3	ug/L		1.32	4.4	7/30/2001	JF
SW-846-8270C	Benzo(k)fluoranthene	<1.6	ug/L		1.58	5.2666	7/30/2001	JF
SW-846-8270C	Benzyl alcohol	<1.8	ug/L		1.8	6	7/30/2001	JF
SW-846-8270C	bis(2-Chloroethoxy)methane	<1.5	ug/L		1.48	4.9333	7/30/2001	JF
SW-846-8270C	bis(2-Chloroethyl)ether	<1.9	ug/L		1.9	6.3333	7/30/2001	JF
SW-846-8270C	bis(2-Chloroisopropyl)ether	<1.5	ug/L		1.54	5.1333	7/30/2001	JF
SW-846-8270C	bis(2-Ethylhexyl)phthalate	<4.6	ug/L		4.6	15.333	7/30/2001	JF
SW-846-8270C	Butylbenzylphthalate	<1.8	ug/L		1.8	6	7/30/2001	JF
SW-846-8270C	Chrysene	<1.6	ug/L		1.56	5.2	7/30/2001	JF
SW-846-8270C	Di-n-butylphthalate	<1.5	ug/L		1.52	5.0666	7/30/2001	JF
SW-846-8270C	Di-n-octylphthalate	<1.5	ug/L		1.52	5.0666	7/30/2001	JF
SW-846-8270C	Dibenz(a,j)acridine	<1.4	ug/L		1.44	4.8	7/30/2001	JF
SW-846-8270C	Dibenzo(a,h)anthracene	<1.5	ug/L		1.5	5	7/30/2001	JF
SW-846-8270C	Dibenzofuran	<1.3	ug/L		1.28	4.2666	7/30/2001	JF
SW-846-8270C	Diethylphthalate	<1.3	ug/L		1.26	4.2	7/30/2001	JF
SW-846-8270C	Dimethylphthalate	<1.5	ug/L		1.52	5.0666	7/30/2001	JF
SW-846-8270C	Diphenylamine	<2.6	ug/L		2.6	8.6666	7/30/2001	JF
SW-846-8270C	Ethyl methanesulfonate	<1.7	ug/L		1.7	5.6666	7/30/2001	JF
SW-846-8270C	Fluoranthene	<2.0	ug/L		1.98	6.6	7/30/2001	JF
SW-846-8270C	Fluorene	<1.2	ug/L		1.2	4	7/30/2001	JF
SW-846-8270C	Hexachlorobenzene	<1.6	ug/L		1.58	5.2666	7/30/2001	JF
SW-846-8270C	Hexachlorobutadiene	<7.6	ug/L		7.6	25.333	7/30/2001	JF
SW-846-8270C	Hexachlorocyclopentadiene	<5.4	ug/L		5.4	18	7/30/2001	JF
SW-846-8270C	Hexachloroethane	<5.2	ug/L		5.2	17.333	7/30/2001	JF
SW-846-8270C	Indeno(1,2,3-cd)pyrene	<1.4	ug/L		1.44	4.8	7/30/2001	JF
SW-846-8270C	Isophorone	<1.4	ug/L		1.42	4.7333	7/30/2001	JF
SW-846-8270C	Methyl methanesulfonate	<1.1	ug/L		1.06	3.5333	7/30/2001	JF
SW-846-8270C	n-Nitrosodi-n-butylamine	<1.5	ug/L		1.46	4.8666	7/30/2001	JF
SW-846-8270C	n-Nitrosodi-n-propylamine	<1.4	ug/L		1.4	4.6666	7/30/2001	JF
SW-846-8270C	n-Nitrosodimethylamine	<1.9	ug/L		1.88	6.2666	7/30/2001	JF
SW-846-8270C	n-Nitrosodiphenylamine	<2.6	ug/L		2.6	8.6666	7/30/2001	JF
SW-846-8270C	n-Nitrosopiperidine	<1.4	ug/L		1.4	4.6666	7/30/2001	JF
SW-846-8270C	Naphthalene	<1.4	ug/L		1.38	4.6	7/30/2001	JF
SW-846-8270C	Nitrobenzene	<1.4	ug/L		1.42	4.7333	7/30/2001	JF
SW-846-8270C	Nitrobenzene-d5 - Surrogate	48	% Rec		0	0	7/30/2001	JF
SW-846-8270C	p-Dimethylaminoazobenzene	<2.0	ug/L		2	6.6666	7/30/2001	JF
SW-846-8270C	Pentachlorobenzene	<2.2	ug/L		2.2	7.3333	7/30/2001	JF
SW-846-8270C	Pentachloronitrobenzene	<1.5	ug/L		1.54	5.1333	7/30/2001	JF
SW-846-8270C	Pentachlorophenol	<1.7	ug/L		1.72	5.7333	7/30/2001	JF
SW-846-8270C	Phenacetin	<1.8	ug/L		1.78	5.9333	7/30/2001	JF

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Method	Parameter Name	Result	Units	Flag	MDL	PQL	Anl Date	Analyst
Lab No	Collect Date	Sample ID						
SW-846-8270C	Phenanthrene	<1.4	ug/L		1.42	4.7333	7/30/2001	JF
SW-846-8270C	Phenol	<1.1	ug/L		1.1	3.6666	7/30/2001	JF
SW-846-8270C	Phenol-d5 - Surrogate	54	% Rec		0	0	7/30/2001	JF
SW-846-8270C	Pronamide	<1.7	ug/L		1.68	5.6	7/30/2001	JF
SW-846-8270C	Pyrene	<2.2	ug/L		2.2	7.3333	7/30/2001	JF
SW-846-8270C	Pyridine	<3.0	ug/L		3	10	7/30/2001	JF
SW-846-8270C	Terphenyl-d14 - Surrogate	52	% Rec		0	0	7/30/2001	JF
01REL010557	6/27/2001	DPZ-1						
EPA 1613	Dioxin Analysis	See Attached			0	0	7/25/2001	PJK
SW-846-8021B	1,2,4-Trimethylbenzene	34	ug/L		4.6	15.333	7/10/2001	CRW
SW-846-8021B	1,3,5-Trimethylbenzene	31	ug/L		4.2	14	7/10/2001	CRW
SW-846-8021B	Benzene	9.7	ug/L	13	4.2	14	7/10/2001	CRW
SW-846-8021B	Ethylbenzene	29	ug/L		4.6	15.333	7/10/2001	CRW
SW-846-8021B	Fluorobenzene-Surrogate	100	% Rec		0	0	7/10/2001	CRW
SW-846-8021B	Methyl-tertiary-butyl ether	<1.8	ug/L		1.82	6.0666	7/10/2001	CRW
SW-846-8021B	Toluene	25	ug/L		4.4	14.666	7/10/2001	CRW
SW-846-8021B	Xylenes-Total	141	ug/L		8.8	29.333	7/10/2001	CRW
SW-846-8270C	Extraction Date	Complete			0	0	7/03/2001	JF
SW-846-8270C	1,2,4,5-Tetrachlorobenzene	<3.0	ug/L		3	10	7/30/2001	JF
SW-846-8270C	1,2,4-Trichlorobenzene	<3.6	ug/L		3.6	12	7/30/2001	JF
SW-846-8270C	1,2-Dichlorobenzene	<3.2	ug/L		3.2	10.666	7/30/2001	JF
SW-846-8270C	1,2-Diphenylhydrazine	<1.3	ug/L		1.34	4.4666	7/30/2001	JF
SW-846-8270C	1,3-Dichlorobenzene	<3.6	ug/L		3.6	12	7/30/2001	JF
SW-846-8270C	1,4-Dichlorobenzene	<3.4	ug/L		3.4	11.333	7/30/2001	JF
SW-846-8270C	1-Chloronaphthalene	<1.8	ug/L		1.84	6.1333	7/30/2001	JF
SW-846-8270C	1-Naphthylamine	<2.2	ug/L		2.2	7.3333	7/30/2001	JF
SW-846-8270C	2,3,4,6-Tetrachlorophenol	44	ug/L		1.36	4.5333	7/30/2001	JF
SW-846-8270C	2,4,5-Trichlorophenol	<1.6	ug/L		1.64	5.4666	7/30/2001	JF
SW-846-8270C	2,4,6-Tribromophenol - Surrogate	62	% Rec		0	0	7/30/2001	JF
SW-846-8270C	2,4,6-Trichlorophenol	<1.6	ug/L		1.62	5.4	7/30/2001	JF
SW-846-8270C	2,4-Dichlorophenol	<1.5	ug/L		1.54	5.1333	7/30/2001	JF
SW-846-8270C	2,4-Dimethylphenol	<2.8	ug/L		2.8	9.3333	7/30/2001	JF
SW-846-8270C	2,4-Dinitrophenol	<1.8	ug/L		1.78	5.9333	7/30/2001	JF
SW-846-8270C	2,4-Dinitrotoluene	<1.6	ug/L		1.58	5.2666	7/30/2001	JF
SW-846-8270C	2,6-Dichlorophenol	<1.5	ug/L		1.52	5.0666	7/30/2001	JF
SW-846-8270C	2,6-Dinitrotoluene	<1.6	ug/L		1.6	5.3333	7/30/2001	JF
SW-846-8270C	2-Chloronaphthalene	<1.6	ug/L		1.56	5.2	7/30/2001	JF
SW-846-8270C	2-Chlorophenol	<1.2	ug/L		1.2	4	7/30/2001	JF
SW-846-8270C	2-Fluorobiphenyl - Surrogate	57	% Rec		0	0	7/30/2001	JF
SW-846-8270C	2-Fluorophenol - Surrogate	43	% Rec		0	0	7/30/2001	JF
SW-846-8270C	2-Methylnaphthalene	3.7	ug/L	13	1.6	5.3333	7/30/2001	JF
SW-846-8270C	2-Methylphenol	<1.7	ug/L		1.7	5.6666	7/30/2001	JF
SW-846-8270C	2-Naphthylamine	<2.2	ug/L		2.2	7.3333	7/30/2001	JF
SW-846-8270C	2-Nitroaniline	<1.6	ug/L		1.64	5.4666	7/30/2001	JF
SW-846-8270C	2-Nitrophenol	<1.3	ug/L		1.28	4.2666	7/30/2001	JF
SW-846-8270C	2-Picoline	<3.0	ug/L		3	10	7/30/2001	JF

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Green Bay, WI 54301-2878

Project Number: 13551002

Project Name: WEISENBERGER TIE & LUMBER

Client ID: L14

Chain: 85968

Report Date: 8/23/2001

Method	Parameter Name	Result	Units	Flag	MDL	PQL	Anl Date	Analyst
Lab No.	Collect Date	Sample ID						
SW-846-8270C	3 & 4-Methylphenol	<1.8	ug/L		1.76	5.8666	7/30/2001	JF
SW-846-8270C	3,3'-Dichlorobenzidine	<0.84	ug/L		0.84	2.8	7/30/2001	JF
SW-846-8270C	3-Methylcholanthrene	<1.9	ug/L		1.94	6.4666	7/30/2001	JF
SW-846-8270C	3-Nitroaniline	<2.2	ug/L		2.2	7.3333	7/30/2001	JF
SW-846-8270C	4,6-Dinitro-2-methylphenol	<2.6	ug/L		2.6	8.6666	7/30/2001	JF
SW-846-8270C	4-Aminobiphenyl	<1.5	ug/L		1.5	5	7/30/2001	JF
SW-846-8270C	4-Bromophenyl phenyl ether	<1.2	ug/L		1.2	4	7/30/2001	JF
SW-846-8270C	4-Chloro-3-methylphenol	<1.7	ug/L		1.7	5.6666	7/30/2001	JF
SW-846-8270C	4-Chloroaniline	<1.3	ug/L		1.32	4.4	7/30/2001	JF
SW-846-8270C	4-Chlorophenyl phenyl ether	<1.2	ug/L		1.24	4.1333	7/30/2001	JF
SW-846-8270C	4-Nitroaniline	<1.3	ug/L		1.32	4.4	7/30/2001	JF
SW-846-8270C	4-Nitrophenol	<1.9	ug/L		1.86	6.2	7/30/2001	JF
SW-846-8270C	7,12-Dimethylbenz(a)anthracene	<2.0	ug/L		2	6.6666	7/30/2001	JF
SW-846-8270C	Acenaphthene	<1.1	ug/L		1.12	3.7333	7/30/2001	JF
SW-846-8270C	Acenaphthylene	<1.5	ug/L		1.46	4.8666	7/30/2001	JF
SW-846-8270C	Acetophenone	1.8	ug/L	13	1.6	5.3333	7/30/2001	JF
SW-846-8270C	Aniline	<1.7	ug/L		1.74	5.8	7/30/2001	JF
SW-846-8270C	Anthracene	<1.3	ug/L		1.3	4.3333	7/30/2001	JF
SW-846-8270C	Benzidine	<2.8	ug/L		2.8	9.3333	7/30/2001	JF
SW-846-8270C	Benzo(a)anthracene	<1.4	ug/L		1.42	4.7333	7/30/2001	JF
SW-846-8270C	Benzo(a)pyrene	<1.1	ug/L		1.1	3.6666	7/30/2001	JF
SW-846-8270C	Benzo(b)fluoranthene	<1.3	ug/L		1.28	4.2666	7/30/2001	JF
SW-846-8270C	Benzo(g,h,i)perylene	<1.3	ug/L		1.32	4.4	7/30/2001	JF
SW-846-8270C	Benzo(k)fluoranthene	<1.6	ug/L		1.58	5.2666	7/30/2001	JF
SW-846-8270C	Benzyl alcohol	<1.8	ug/L		1.8	6	7/30/2001	JF
SW-846-8270C	bis(2-Chloroethoxy)methane	<1.5	ug/L		1.48	4.9333	7/30/2001	JF
SW-846-8270C	bis(2-Chloroethyl)ether	<1.9	ug/L		1.9	6.3333	7/30/2001	JF
SW-846-8270C	bis(2-Chloroisopropyl)ether	<1.5	ug/L		1.54	5.1333	7/30/2001	JF
SW-846-8270C	bis(2-Ethylhexyl)phthalate	<4.6	ug/L		4.6	15.333	7/30/2001	JF
SW-846-8270C	Butylbenzylphthalate	<1.8	ug/L		1.8	6	7/30/2001	JF
SW-846-8270C	Chrysene	<1.6	ug/L		1.56	5.2	7/30/2001	JF
SW-846-8270C	Di-n-butylphthalate	<1.5	ug/L		1.52	5.0666	7/30/2001	JF
SW-846-8270C	Di-n-octylphthalate	<1.5	ug/L		1.52	5.0666	7/30/2001	JF
SW-846-8270C	Dibenz(a,j)acridine	<1.4	ug/L		1.44	4.8	7/30/2001	JF
SW-846-8270C	Dibenzo(a,h)anthracene	<1.5	ug/L		1.5	5	7/30/2001	JF
SW-846-8270C	Dibenzofuran	<1.3	ug/L		1.28	4.2666	7/30/2001	JF
SW-846-8270C	Diethylphthalate	<1.3	ug/L		1.26	4.2	7/30/2001	JF
SW-846-8270C	Dimethylphthalate	<1.5	ug/L		1.52	5.0666	7/30/2001	JF
SW-846-8270C	Diphenylamine	<2.6	ug/L		2.6	8.6666	7/30/2001	JF
SW-846-8270C	Ethyl methanesulfonate	<1.7	ug/L		1.7	5.6666	7/30/2001	JF
SW-846-8270C	Fluoranthene	<2.0	ug/L		1.98	6.6	7/30/2001	JF
SW-846-8270C	Fluorene	<1.2	ug/L		1.2	4	7/30/2001	JF
SW-846-8270C	Hexachlorobenzene	<1.6	ug/L		1.58	5.2666	7/30/2001	JF
SW-846-8270C	Hexachlorobutadiene	<7.6	ug/L		7.6	25.333	7/30/2001	JF
SW-846-8270C	Hexachlorocyclopentadiene	<5.4	ug/L		5.4	18	7/30/2001	JF
SW-846-8270C	Hexachloroethane	<5.2	ug/L		5.2	17.333	7/30/2001	JF

Robert E. Lee & Associates, Inc.
Wisconsin Certification Number: 405043870
Certificate of Analysis Report

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Green Bay, WI 54301-2878

Client ID: L14

Project Number: 13551002

Chain: 85968

Project Name: WEISENBERGER TIE & LUMBER

Report Date: 8/23/2001

Method	Parameter Name	Result	Units	Flag	MDL	PQL	Anl Date	Analyst
Lab No.	Collect Date	Sample ID						
SW-846-8270C	Indeno(1,2,3-cd)pyrene	<1.4	ug/L		1.44	4.8	7/30/2001	JF
SW-846-8270C	Isophorone	5.9	ug/L		1.42	4.7333	7/30/2001	JF
SW-846-8270C	Methyl methanesulfonate	<1.1	ug/L		1.06	3.5333	7/30/2001	JF
SW-846-8270C	n-Nitrosodi-n-butylamine	<1.5	ug/L		1.46	4.8666	7/30/2001	JF
SW-846-8270C	n-Nitrosodi-n-propylamine	<1.4	ug/L		1.4	4.6666	7/30/2001	JF
SW-846-8270C	n-Nitrosodimethylamine	<1.9	ug/L		1.88	6.2666	7/30/2001	JF
SW-846-8270C	n-Nitrosodiphenylamine	<2.6	ug/L		2.6	8.6666	7/30/2001	JF
SW-846-8270C	n-Nitrosopiperidine	<1.4	ug/L		1.4	4.6666	7/30/2001	JF
SW-846-8270C	Naphthalene	2.18	ug/L	13	1.38	4.6	7/30/2001	JF
SW-846-8270C	Nitrobenzene	<1.4	ug/L		1.42	4.7333	7/30/2001	JF
SW-846-8270C	Nitrobenzene-d5 - Surrogate	50	% Rec		0	0	7/30/2001	JF
SW-846-8270C	p-Dimethylaminoazobenzene	<2.0	ug/L		2	6.6666	7/30/2001	JF
SW-846-8270C	Pentachlorobenzene	<2.2	ug/L		2.2	7.3333	7/30/2001	JF
SW-846-8270C	Pentachloronitrobenzene	<1.5	ug/L		1.54	5.1333	7/30/2001	JF
SW-846-8270C	Pentachlorophenol	3120	ug/L		1.72	5.7333	7/30/2001	JF
SW-846-8270C	Phenacetin	<1.8	ug/L		1.78	5.9333	7/30/2001	JF
SW-846-8270C	Phenanthrene	<1.4	ug/L		1.42	4.7333	7/30/2001	JF
SW-846-8270C	Phenol	<1.1	ug/L		1.1	3.6666	7/30/2001	JF
SW-846-8270C	Phenol-d5 - Surrogate	39	% Rec		0	0	7/30/2001	JF
SW-846-8270C	Pronamide	<1.7	ug/L		1.68	5.6	7/30/2001	JF
SW-846-8270C	Pyrene	<2.2	ug/L		2.2	7.3333	7/30/2001	JF
SW-846-8270C	Pyridine	<3.0	ug/L		3	10	7/30/2001	JF
SW-846-8270C	Terphenyl-d14 - Surrogate	50	% Rec		0	0	7/30/2001	JF
01RELO10558	6/27/2001	DPZ-1A						
SW-846-8270C	Extraction Date	Complete			0	0	7/03/2001	JF
SW-846-8270C	1,2,4,5-Tetrachlorobenzene	<3.0	ug/L		3	10	7/30/2001	JF
SW-846-8270C	1,2,4-Trichlorobenzene	<3.6	ug/L		3.6	12	7/30/2001	JF
SW-846-8270C	1,2-Dichlorobenzene	<3.2	ug/L		3.2	10.666	7/30/2001	JF
SW-846-8270C	1,2-Diphenylhydrazine	<1.3	ug/L		1.34	4.4666	7/30/2001	JF
SW-846-8270C	1,3-Dichlorobenzene	<3.6	ug/L		3.6	12	7/30/2001	JF
SW-846-8270C	1,4-Dichlorobenzene	<3.4	ug/L		3.4	11.333	7/30/2001	JF
SW-846-8270C	1-Chloronaphthalene	<1.8	ug/L		1.84	6.1333	7/30/2001	JF
SW-846-8270C	1-Naphthylamine	<2.2	ug/L		2.2	7.3333	7/30/2001	JF
SW-846-8270C	2,3,4,6-Tetrachlorophenol	<1.4	ug/L		1.36	4.5333	7/30/2001	JF
SW-846-8270C	2,4,5-Trichlorophenol	<1.6	ug/L		1.64	5.4666	7/30/2001	JF
SW-846-8270C	2,4,6-Tribromophenol - Surrogate	53	% Rec		0	0	7/30/2001	JF
SW-846-8270C	2,4,6-Trichlorophenol	<1.6	ug/L		1.62	5.4	7/30/2001	JF
SW-846-8270C	2,4-Dichlorophenol	<1.5	ug/L		1.54	5.1333	7/30/2001	JF
SW-846-8270C	2,4-Dimethylphenol	<2.8	ug/L		2.8	9.3333	7/30/2001	JF
SW-846-8270C	2,4-Dinitrophenol	<1.8	ug/L		1.78	5.9333	7/30/2001	JF
SW-846-8270C	2,4-Dinitrotoluene	<1.6	ug/L		1.58	5.2666	7/30/2001	JF
SW-846-8270C	2,6-Dichlorophenol	<1.5	ug/L		1.52	5.0666	7/30/2001	JF
SW-846-8270C	2,6-Dinitrotoluene	<1.6	ug/L		1.6	5.3333	7/30/2001	JF
SW-846-8270C	2-Chloronaphthalene	<1.6	ug/L		1.56	5.2	7/30/2001	JF
SW-846-8270C	2-Chlorophenol	<1.2	ug/L		1.2	4	7/30/2001	JF
SW-846-8270C	2-Fluorobiphenyl - Surrogate	48	% Rec		0	0	7/30/2001	JF

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 Client ID: L14
 Chain: 85968
 Report Date: 8/23/2001

Method	Parameter Name	Result	Units	Flag	MDL	PQL	Anal Date	Analyst
Lab No.	Collect Date	Sample ID						
SW-846-8270C	2-Fluorophenol - Surrogate	32	% Rec		0	0	7/30/2001	JF
SW-846-8270C	2-Methylnaphthalene	<1.6	ug/L		1.6	5.3333	7/30/2001	JF
SW-846-8270C	2-Methylphenol	<1.7	ug/L		1.7	5.6666	7/30/2001	JF
SW-846-8270C	2-Naphthylamine	<2.2	ug/L		2.2	7.3333	7/30/2001	JF
SW-846-8270C	2-Nitroaniline	<1.6	ug/L		1.64	5.4666	7/30/2001	JF
SW-846-8270C	2-Nitrophenol	<1.3	ug/L		1.28	4.2666	7/30/2001	JF
SW-846-8270C	2-Picoline	<3.0	ug/L		3	10	7/30/2001	JF
SW-846-8270C	3 & 4-Methylphenol	<1.8	ug/L		1.76	5.8666	7/30/2001	JF
SW-846-8270C	3,3'-Dichlorobenzidine	<0.84	ug/L		0.84	2.8	7/30/2001	JF
SW-846-8270C	3-Methylcholanthrene	<1.9	ug/L		1.94	6.4666	7/30/2001	JF
SW-846-8270C	3-Nitroaniline	<2.2	ug/L		2.2	7.3333	7/30/2001	JF
SW-846-8270C	4,6-Dinitro-2-methylphenol	<2.6	ug/L		2.6	8.6666	7/30/2001	JF
SW-846-8270C	4-Aminobiphenyl	<1.5	ug/L		1.5	5	7/30/2001	JF
SW-846-8270C	4-Bromophenyl phenyl ether	<1.2	ug/L		1.2	4	7/30/2001	JF
SW-846-8270C	4-Chloro-3-methylphenol	<1.7	ug/L		1.7	5.6666	7/30/2001	JF
SW-846-8270C	4-Chloroaniline	<1.3	ug/L		1.32	4.4	7/30/2001	JF
SW-846-8270C	4-Chlorophenyl phenyl ether	<1.2	ug/L		1.24	4.1333	7/30/2001	JF
SW-846-8270C	4-Nitroaniline	<1.3	ug/L		1.32	4.4	7/30/2001	JF
SW-846-8270C	4-Nitrophenol	<1.9	ug/L		1.86	6.2	7/30/2001	JF
SW-846-8270C	7,12-Dimethylbenz(a)anthracene	<2.0	ug/L		2	6.6666	7/30/2001	JF
SW-846-8270C	Acenaphthene	<1.1	ug/L		1.12	3.7333	7/30/2001	JF
SW-846-8270C	Acenaphthylene	<1.5	ug/L		1.46	4.8666	7/30/2001	JF
SW-846-8270C	Acetophenone	<1.6	ug/L		1.6	5.3333	7/30/2001	JF
SW-846-8270C	Aniline	<1.7	ug/L		1.74	5.8	7/30/2001	JF
SW-846-8270C	Anthracene	<1.3	ug/L		1.3	4.3333	7/30/2001	JF
SW-846-8270C	Benzidine	<2.8	ug/L		2.8	9.3333	7/30/2001	JF
SW-846-8270C	Benzo(a)anthracene	<1.4	ug/L		1.42	4.7333	7/30/2001	JF
SW-846-8270C	Benzo(a)pyrene	<1.1	ug/L		1.1	3.6666	7/30/2001	JF
SW-846-8270C	Benzo(b)fluoranthene	<1.3	ug/L		1.28	4.2666	7/30/2001	JF
SW-846-8270C	Benzo(g,h,i)perylene	<1.3	ug/L		1.32	4.4	7/30/2001	JF
SW-846-8270C	Benzo(k)fluoranthene	<1.6	ug/L		1.58	5.2666	7/30/2001	JF
SW-846-8270C	Benzyl alcohol	<1.8	ug/L		1.8	6	7/30/2001	JF
SW-846-8270C	bis(2-Chloroethoxy)methane	<1.5	ug/L		1.48	4.9333	7/30/2001	JF
SW-846-8270C	bis(2-Chloroethyl)ether	<1.9	ug/L		1.9	6.3333	7/30/2001	JF
SW-846-8270C	bis(2-Chloroisopropyl)ether	<1.5	ug/L		1.54	5.1333	7/30/2001	JF
SW-846-8270C	bis(2-Ethylhexyl)phthalate	<4.6	ug/L		4.6	15.333	7/30/2001	JF
SW-846-8270C	Butylbenzylphthalate	<1.8	ug/L		1.8	6	7/30/2001	JF
SW-846-8270C	Chrysene	<1.6	ug/L		1.56	5.2	7/30/2001	JF
SW-846-8270C	Di-n-butylphthalate	<1.5	ug/L		1.52	5.0666	7/30/2001	JF
SW-846-8270C	Di-n-octylphthalate	<1.5	ug/L		1.52	5.0666	7/30/2001	JF
SW-846-8270C	Dibenz(a,j)acridine	<1.4	ug/L		1.44	4.8	7/30/2001	JF
SW-846-8270C	Dibenzo(a,h)anthracene	<1.5	ug/L		1.5	5	7/30/2001	JF
SW-846-8270C	Dibenzofuran	<1.3	ug/L		1.28	4.2666	7/30/2001	JF
SW-846-8270C	Diethylphthalate	<1.3	ug/L		1.26	4.2	7/30/2001	JF
SW-846-8270C	Dimethylphthalate	<1.5	ug/L		1.52	5.0666	7/30/2001	JF
SW-846-8270C	Diphenylamine	<2.6	ug/L		2.6	8.6666	7/30/2001	JF

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Client ID: L14
Chain: 85968
Report Date: 8/23/2001

Method	Parameter Name	Result	Units	Flag	MDL	PQL	Anl Date	Analyst
Lab No.	Collect Date	Sample ID						
SW-846-8270C	Ethyl methanesulfonate	<1.7	ug/L		1.7	5.6666	7/30/2001	JF
SW-846-8270C	Fluoranthene	<2.0	ug/L		1.98	6.6	7/30/2001	JF
SW-846-8270C	Fluorene	<1.2	ug/L		1.2	4	7/30/2001	JF
SW-846-8270C	Hexachlorobenzene	<1.6	ug/L		1.58	5.2666	7/30/2001	JF
SW-846-8270C	Hexachlorobutadiene	<7.6	ug/L		7.6	25.333	7/30/2001	JF
SW-846-8270C	Hexachlorocyclopentadiene	<5.4	ug/L		5.4	18	7/30/2001	JF
SW-846-8270C	Hexachloroethane	<5.2	ug/L		5.2	17.333	7/30/2001	JF
SW-846-8270C	Indeno(1,2,3-cd)pyrene	<1.4	ug/L		1.44	4.8	7/30/2001	JF
SW-846-8270C	Isophorone	<1.4	ug/L		1.42	4.7333	7/30/2001	JF
SW-846-8270C	Methyl methanesulfonate	<1.1	ug/L		1.06	3.5333	7/30/2001	JF
SW-846-8270C	n-Nitrosodi-n-butylamine	<1.5	ug/L		1.46	4.8666	7/30/2001	JF
SW-846-8270C	n-Nitrosodi-n-propylamine	<1.4	ug/L		1.4	4.6666	7/30/2001	JF
SW-846-8270C	n-Nitrosodimethylamine	<1.9	ug/L		1.88	6.2666	7/30/2001	JF
SW-846-8270C	n-Nitrosodiphenylamine	<2.6	ug/L		2.6	8.6666	7/30/2001	JF
SW-846-8270C	n-Nitrosopiperidine	<1.4	ug/L		1.4	4.6666	7/30/2001	JF
SW-846-8270C	Naphthalene	<1.4	ug/L		1.38	4.6	7/30/2001	JF
SW-846-8270C	Nitrobenzene	<1.4	ug/L		1.42	4.7333	7/30/2001	JF
SW-846-8270C	Nitrobenzene-d5 - Surrogate	39	% Rec		0	0	7/30/2001	JF
SW-846-8270C	p-Dimethylaminoazobenzene	<2.0	ug/L		2	6.6666	7/30/2001	JF
SW-846-8270C	Pentachlorobenzene	<2.2	ug/L		2.2	7.3333	7/30/2001	JF
SW-846-8270C	Pentachloronitrobenzene	<1.5	ug/L		1.54	5.1333	7/30/2001	JF
SW-846-8270C	Pentachlorophenol	<1.7	ug/L		1.72	5.7333	7/30/2001	JF
SW-846-8270C	Phenacetin	<1.8	ug/L		1.78	5.9333	7/30/2001	JF
SW-846-8270C	Phenanthrene	<1.4	ug/L		1.42	4.7333	7/30/2001	JF
SW-846-8270C	Phenol	<1.1	ug/L		1.1	3.6666	7/30/2001	JF
SW-846-8270C	Phenol-d5 - Surrogate	30	% Rec		0	0	7/30/2001	JF
SW-846-8270C	Pronamide	<1.7	ug/L		1.68	5.6	7/30/2001	JF
SW-846-8270C	Pyrene	<2.2	ug/L		2.2	7.3333	7/30/2001	JF
SW-846-8270C	Pyridine	<3.0	ug/L		3	10	7/30/2001	JF
SW-846-8270C	Terphenyl-d14 - Surrogate	52	% Rec		0	0	7/30/2001	JF
01REL010559	6/27/2001	DPZ-2						
SW-846-8270C	Extraction Date	Complete			0	0	7/03/2001	JF
SW-846-8270C	1,2,4,5-Tetrachlorobenzene	<3.0	ug/L		3	10	7/30/2001	JF
SW-846-8270C	1,2,4-Trichlorobenzene	<3.6	ug/L		3.6	12	7/30/2001	JF
SW-846-8270C	1,2-Dichlorobenzene	<3.2	ug/L		3.2	10.666	7/30/2001	JF
SW-846-8270C	1,2-Diphenylhydrazine	<1.3	ug/L		1.34	4.4666	7/30/2001	JF
SW-846-8270C	1,3-Dichlorobenzene	<3.6	ug/L		3.6	12	7/30/2001	JF
SW-846-8270C	1,4-Dichlorobenzene	<3.4	ug/L		3.4	11.333	7/30/2001	JF
SW-846-8270C	1-Chloronaphthalene	<1.8	ug/L		1.84	6.1333	7/30/2001	JF
SW-846-8270C	1-Naphthylamine	<2.2	ug/L		2.2	7.3333	7/30/2001	JF
SW-846-8270C	2,3,4,6-Tetrachlorophenol	<1.4	ug/L		1.36	4.5333	7/30/2001	JF
SW-846-8270C	2,4,5-Trichlorophenol	<1.6	ug/L		1.64	5.4666	7/30/2001	JF
SW-846-8270C	2,4,6-Tribromophenol - Surrogate	64	% Rec		0	0	7/30/2001	JF
SW-846-8270C	2,4,6-Trichlorophenol	<1.6	ug/L		1.62	5.4	7/30/2001	JF
SW-846-8270C	2,4-Dichlorophenol	<1.5	ug/L		1.54	5.1333	7/30/2001	JF
SW-846-8270C	2,4-Dimethylphenol	<2.8	ug/L		2.8	9.3333	7/30/2001	JF

Robert E Lee & Associates, Inc
 2825 S Webster Ave

Green Bay, WI 54301-2878

Project Number: 13551002

Project Name: WEISENBERGER TIE & LUMBER

Attn: Jim Caine

Phone: (920)336-6338

Fax: (920)336-9141

Client ID: L14

Chain: 85968

Report Date: 8/23/2001

Method	Parameter Name	Result	Units	Flag	MDL	PQL	Anl Date	Analyst
Lab No.	Collect Date	Sample ID						
SW-846-8270C	2,4-Dinitrophenol	<1.8	ug/L		1.78	5.9333	7/30/2001	JF
SW-846-8270C	2,4-Dinitrotoluene	<1.6	ug/L		1.58	5.2666	7/30/2001	JF
SW-846-8270C	2,6-Dichlorophenol	<1.5	ug/L		1.52	5.0666	7/30/2001	JF
SW-846-8270C	2,6-Dinitrotoluene	<1.6	ug/L		1.6	5.3333	7/30/2001	JF
SW-846-8270C	2-Chloronaphthalene	<1.6	ug/L		1.56	5.2	7/30/2001	JF
SW-846-8270C	2-Chlorophenol	<1.2	ug/L		1.2	4	7/30/2001	JF
SW-846-8270C	2-Fluorobiphenyl - Surrogate	55	% Rec		0	0	7/30/2001	JF
SW-846-8270C	2-Fluorophenol - Surrogate	34	% Rec		0	0	7/30/2001	JF
SW-846-8270C	2-Methylnaphthalene	<1.6	ug/L		1.6	5.3333	7/30/2001	JF
SW-846-8270C	2-Methylphenol	<1.7	ug/L		1.7	5.6666	7/30/2001	JF
SW-846-8270C	2-Naphthylamine	<2.2	ug/L		2.2	7.3333	7/30/2001	JF
SW-846-8270C	2-Nitroaniline	<1.6	ug/L		1.64	5.4666	7/30/2001	JF
SW-846-8270C	2-Nitrophenol	<1.3	ug/L		1.28	4.2666	7/30/2001	JF
SW-846-8270C	2-Picoline	<3.0	ug/L		3	10	7/30/2001	JF
SW-846-8270C	3 & 4-Methylphenol	<1.8	ug/L		1.76	5.8666	7/30/2001	JF
SW-846-8270C	3,3'-Dichlorobenzidine	<0.84	ug/L		0.84	2.8	7/30/2001	JF
SW-846-8270C	3-Methylcholanthrene	<1.9	ug/L		1.94	6.4666	7/30/2001	JF
SW-846-8270C	3-Nitroaniline	<2.2	ug/L		2.2	7.3333	7/30/2001	JF
SW-846-8270C	4,6-Dinitro-2-methylphenol	<2.6	ug/L		2.6	8.6666	7/30/2001	JF
SW-846-8270C	4-Aminobiphenyl	<1.5	ug/L		1.5	5	7/30/2001	JF
SW-846-8270C	4-Bromophenyl phenyl ether	<1.2	ug/L		1.2	4	7/30/2001	JF
SW-846-8270C	4-Chloro-3-methylphenol	<1.7	ug/L		1.7	5.6666	7/30/2001	JF
SW-846-8270C	4-Chloroaniline	<1.3	ug/L		1.32	4.4	7/30/2001	JF
SW-846-8270C	4-Chlorophenyl phenyl ether	<1.2	ug/L		1.24	4.1333	7/30/2001	JF
SW-846-8270C	4-Nitroaniline	<1.3	ug/L		1.32	4.4	7/30/2001	JF
SW-846-8270C	4-Nitrophenol	<1.9	ug/L		1.86	6.2	7/30/2001	JF
SW-846-8270C	7,12-Dimethylbenz(a)anthracene	<2.0	ug/L		2	6.6666	7/30/2001	JF
SW-846-8270C	Acenaphthene	<1.1	ug/L		1.12	3.7333	7/30/2001	JF
SW-846-8270C	Acenaphthylene	<1.5	ug/L		1.46	4.8666	7/30/2001	JF
SW-846-8270C	Acetophenone	<1.6	ug/L		1.6	5.3333	7/30/2001	JF
SW-846-8270C	Aniline	<1.7	ug/L		1.74	5.8	7/30/2001	JF
SW-846-8270C	Anthracene	<1.3	ug/L		1.3	4.3333	7/30/2001	JF
SW-846-8270C	Benzidine	<2.8	ug/L		2.8	9.3333	7/30/2001	JF
SW-846-8270C	Benzo(a)anthracene	<1.4	ug/L		1.42	4.7333	7/30/2001	JF
SW-846-8270C	Benzo(a)pyrene	<1.1	ug/L		1.1	3.6666	7/30/2001	JF
SW-846-8270C	Benzo(b)fluoranthene	<1.3	ug/L		1.28	4.2666	7/30/2001	JF
SW-846-8270C	Benzo(g,h,i)perylene	<1.3	ug/L		1.32	4.4	7/30/2001	JF
SW-846-8270C	Benzo(k)fluoranthene	<1.6	ug/L		1.58	5.2666	7/30/2001	JF
SW-846-8270C	Benzyl alcohol	<1.8	ug/L		1.8	6	7/30/2001	JF
SW-846-8270C	bis(2-Chloroethoxy)methane	<1.5	ug/L		1.48	4.9333	7/30/2001	JF
SW-846-8270C	bis(2-Chloroethyl)ether	<1.9	ug/L		1.9	6.3333	7/30/2001	JF
SW-846-8270C	bis(2-Chloroisopropyl)ether	<1.5	ug/L		1.54	5.1333	7/30/2001	JF
SW-846-8270C	bis(2-Ethylhexyl)phthalate	<4.6	ug/L		4.6	15.333	7/30/2001	JF
SW-846-8270C	Butylbenzylphthalate	<1.8	ug/L		1.8	6	7/30/2001	JF
SW-846-8270C	Chrysene	<1.6	ug/L		1.56	5.2	7/30/2001	JF
SW-846-8270C	Di-n-butylphthalate	<1.5	ug/L		1.52	5.0666	7/30/2001	JF

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 Client ID: L14
 Chain: 85968
 Report Date: 8/23/2001

Green Bay, WI 54301-2878
 Project Number: 13551002
 Project Name: WEISENBERGER TIE & LUMBER

Method	Parameter Name	Result	Units	Flag	MDL	PQL	Anl Date	Analyst
Lab No	Collect Date	Sample ID						
SW-846-8270C	Di-n-octylphthalate	<1.5	ug/L		1.52	5.0666	7/30/2001	JF
SW-846-8270C	Dibenz(a,j)acridine	<1.4	ug/L		1.44	4.8	7/30/2001	JF
SW-846-8270C	Dibenzo(a,h)anthracene	<1.5	ug/L		1.5	5	7/30/2001	JF
SW-846-8270C	Dibenzofuran	<1.3	ug/L		1.28	4.2666	7/30/2001	JF
SW-846-8270C	Diethylphthalate	<1.3	ug/L		1.26	4.2	7/30/2001	JF
SW-846-8270C	Dlmethylphthalate	<1.5	ug/L		1.52	5.0666	7/30/2001	JF
SW-846-8270C	Diphenylamine	<2.6	ug/L		2.6	8.6666	7/30/2001	JF
SW-846-8270C	Ethyl methanesulfonate	<1.7	ug/L		1.7	5.6666	7/30/2001	JF
SW-846-8270C	Fluoranthene	<2.0	ug/L		1.98	6.6	7/30/2001	JF
SW-846-8270C	Fluorene	<1.2	ug/L		1.2	4	7/30/2001	JF
SW-846-8270C	Hexachlorobenzene	<1.6	ug/L		1.58	5.2666	7/30/2001	JF
SW-846-8270C	Hexachlorobutadiene	<7.6	ug/L		7.6	25.333	7/30/2001	JF
SW-846-8270C	Hexachlorocyclopentadiene	<5.4	ug/L		5.4	18	7/30/2001	JF
SW-846-8270C	Hexachloroethane	<5.2	ug/L		5.2	17.333	7/30/2001	JF
SW-846-8270C	Indeno(1,2,3-cd)pyrene	<1.4	ug/L		1.44	4.8	7/30/2001	JF
SW-846-8270C	Isophorone	<1.4	ug/L		1.42	4.7333	7/30/2001	JF
SW-846-8270C	Methyl methanesulfonate	<1.1	ug/L		1.06	3.5333	7/30/2001	JF
SW-846-8270C	n-Nitrosodi-n-butylamine	<1.5	ug/L		1.46	4.8666	7/30/2001	JF
SW-846-8270C	n-Nitrosodi-n-propylamine	<1.4	ug/L		1.4	4.6666	7/30/2001	JF
SW-846-8270C	n-Nitrosodimethylamine	<1.9	ug/L		1.88	6.2666	7/30/2001	JF
SW-846-8270C	n-Nitrosodiphenylamine	<2.6	ug/L		2.6	8.6666	7/30/2001	JF
SW-846-8270C	n-Nitrosopiperidine	<1.4	ug/L		1.4	4.6666	7/30/2001	JF
SW-846-8270C	Naphthalene	<1.4	ug/L		1.38	4.6	7/30/2001	JF
SW-846-8270C	Nitrobenzene	<1.4	ug/L		1.42	4.7333	7/30/2001	JF
SW-846-8270C	Nitrobenzene-d5 - Surrogate	43	% Rec		0	0	7/30/2001	JF
SW-846-8270C	p-Dimethylaminoazobenzene	<2.0	ug/L		2	6.6666	7/30/2001	JF
SW-846-8270C	Pentachlorobenzene	<2.2	ug/L		2.2	7.3333	7/30/2001	JF
SW-846-8270C	Pentachloronitrobenzene	<1.5	ug/L		1.54	5.1333	7/30/2001	JF
SW-846-8270C	Pentachlorophenol	1.9	ug/L	13	1.72	5.7333	7/30/2001	JF
SW-846-8270C	Phenacetin	<1.8	ug/L		1.78	5.9333	7/30/2001	JF
SW-846-8270C	Phenanthrene	<1.4	ug/L		1.42	4.7333	7/30/2001	JF
SW-846-8270C	Phenol	<1.1	ug/L		1.1	3.6666	7/30/2001	JF
SW-846-8270C	Phenol-d5 - Surrogate	31	% Rec		0	0	7/30/2001	JF
SW-846-8270C	Pronamide	<1.7	ug/L		1.68	5.6	7/30/2001	JF
SW-846-8270C	Pyrene	<2.2	ug/L		2.2	7.3333	7/30/2001	JF
SW-846-8270C	Pyridine	<3.0	ug/L		3	10	7/30/2001	JF
SW-846-8270C	Terphenyl-d14 - Surrogate	55	% Rec		0	0	7/30/2001	JF
01REL010560 6/27/2001 DPZ-3								
EPA 1613	Dioxin Analysis	See Attached			0	0	7/25/2001	PJK
SW-846-8270C	Extraction Date	Complete			0	0	7/03/2001	JF
SW-846-8270C	1,2,4,5-Tetrachlorobenzene	<3.0	ug/L		3	10	7/30/2001	JF
SW-846-8270C	1,2,4-Trichlorobenzene	<3.6	ug/L		3.6	12	7/30/2001	JF
SW-846-8270C	1,2-Dichlorobenzene	<3.2	ug/L		3.2	10.666	7/30/2001	JF
SW-846-8270C	1,2-Diphenylhydrazine	<1.3	ug/L		1.34	4.4666	7/30/2001	JF
SW-846-8270C	1,3-Dichlorobenzene	<3.6	ug/L		3.6	12	7/30/2001	JF
SW-846-8270C	1,4-Dichlorobenzene	<3.4	ug/L		3.4	11.333	7/30/2001	JF

Robert E Lee & Associates, Inc
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Project Number: 13551002

Project Name: WEISENBERGER TIE & LUMBER

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Client ID: L14

Chain: 85968

Report Date: 8/23/2001

Method	Parameter Name	Result	Units	Flag	MDL	PQL	Anl Date	Analyst
Lab No.	Collect Date	Sample ID						
SW-846-8270C		1-Chloronaphthalene	<1.8	ug/L	1.84	6.1333	7/30/2001	JF
SW-846-8270C		1-Naphthylamine	<2.2	ug/L	2.2	7.3333	7/30/2001	JF
SW-846-8270C		2,3,4,6-Tetrachlorophenol	<1.4	ug/L	1.36	4.5333	7/30/2001	JF
SW-846-8270C		2,4,5-Trichlorophenol	<1.6	ug/L	1.64	5.4666	7/30/2001	JF
SW-846-8270C		2,4,6-Tribromophenol - Surrogate	71	% Rec	0	0	7/30/2001	JF
SW-846-8270C		2,4,6-Trichlorophenol	<1.6	ug/L	1.62	5.4	7/30/2001	JF
SW-846-8270C		2,4-Dichlorophenol	<1.5	ug/L	1.54	5.1333	7/30/2001	JF
SW-846-8270C		2,4-Dimethylphenol	<2.8	ug/L	2.8	9.3333	7/30/2001	JF
SW-846-8270C		2,4-Dinitrophenol	<1.8	ug/L	1.78	5.9333	7/30/2001	JF
SW-846-8270C		2,4-Dinitrotoluene	<1.6	ug/L	1.58	5.2666	7/30/2001	JF
SW-846-8270C		2,6-Dichlorophenol	<1.5	ug/L	1.52	5.0666	7/30/2001	JF
SW-846-8270C		2,6-Dinitrotoluene	<1.6	ug/L	1.6	5.3333	7/30/2001	JF
SW-846-8270C		2-Chloronaphthalene	<1.6	ug/L	1.56	5.2	7/30/2001	JF
SW-846-8270C		2-Chlorophenol	<1.2	ug/L	1.2	4	7/30/2001	JF
SW-846-8270C		2-Fluorobiphenyl - Surrogate	59	% Rec	0	0	7/30/2001	JF
SW-846-8270C		2-Fluorophenol - Surrogate	45	% Rec	0	0	7/30/2001	JF
SW-846-8270C		2-Methylnaphthalene	<1.6	ug/L	1.6	5.3333	7/30/2001	JF
SW-846-8270C		2-Methylphenol	<1.7	ug/L	1.7	5.6666	7/30/2001	JF
SW-846-8270C		2-Naphthylamine	<2.2	ug/L	2.2	7.3333	7/30/2001	JF
SW-846-8270C		2-Nitroaniline	<1.6	ug/L	1.64	5.4666	7/30/2001	JF
SW-846-8270C		2-Nitrophenol	<1.3	ug/L	1.28	4.2666	7/30/2001	JF
SW-846-8270C		2-Picoline	<3.0	ug/L	3	10	7/30/2001	JF
SW-846-8270C		3 & 4-Methylphenol	<1.8	ug/L	1.76	5.8666	7/30/2001	JF
SW-846-8270C		3,3'-Dichlorobenzidine	<0.84	ug/L	0.84	2.8	7/30/2001	JF
SW-846-8270C		3-Methylcholanthrene	<1.9	ug/L	1.94	6.4666	7/30/2001	JF
SW-846-8270C		3-Nitroaniline	<2.2	ug/L	2.2	7.3333	7/30/2001	JF
SW-846-8270C		4,6-Dinitro-2-methylphenol	<2.6	ug/L	2.6	8.6666	7/30/2001	JF
SW-846-8270C		4-Aminobiphenyl	<1.5	ug/L	1.5	5	7/30/2001	JF
SW-846-8270C		4-Bromophenyl phenyl ether	<1.2	ug/L	1.2	4	7/30/2001	JF
SW-846-8270C		4-Chloro-3-methylphenol	<1.7	ug/L	1.7	5.6666	7/30/2001	JF
SW-846-8270C		4-Chloroaniline	<1.3	ug/L	1.32	4.4	7/30/2001	JF
SW-846-8270C		4-Chlorophenyl phenyl ether	<1.2	ug/L	1.24	4.1333	7/30/2001	JF
SW-846-8270C		4-Nitroaniline	<1.3	ug/L	1.32	4.4	7/30/2001	JF
SW-846-8270C		4-Nitrophenol	<1.9	ug/L	1.86	6.2	7/30/2001	JF
SW-846-8270C		7,12-Dimethylbenz(a)anthracene	<2.0	ug/L	2	6.6666	7/30/2001	JF
SW-846-8270C		Acenaphthene	<1.1	ug/L	1.12	3.7333	7/30/2001	JF
SW-846-8270C		Acenaphthylene	<1.5	ug/L	1.46	4.8666	7/30/2001	JF
SW-846-8270C		Acetophenone	<1.6	ug/L	1.6	5.3333	7/30/2001	JF
SW-846-8270C		Aniline	<1.7	ug/L	1.74	5.8	7/30/2001	JF
SW-846-8270C		Anthracene	<1.3	ug/L	1.3	4.3333	7/30/2001	JF
SW-846-8270C		Benzidine	<2.8	ug/L	2.8	9.3333	7/30/2001	JF
SW-846-8270C		Benzo(a)anthracene	<1.4	ug/L	1.42	4.7333	7/30/2001	JF
SW-846-8270C		Benzo(a)pyrene	<1.1	ug/L	1.1	3.6666	7/30/2001	JF
SW-846-8270C		Benzo(b)fluoranthene	<1.3	ug/L	1.28	4.2666	7/30/2001	JF
SW-846-8270C		Benzo(g,h,i)perylene	<1.3	ug/L	1.32	4.4	7/30/2001	JF
SW-846-8270C		Benzo(k)fluoranthene	<1.6	ug/L	1.58	5.2666	7/30/2001	JF

Robert E. Lee & Associates, Inc.
Wisconsin Certification Number: 405043870
Certificate of Analysis Report

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Method	Parameter Name	Result	Units	Flag	MDL	PQL	Anl Date	Analyst
Lab No.	Collect Date	Sample ID						
SW-846-8270C	Benzyl alcohol	<1.8	ug/L		1.8	6	7/30/2001	JF
SW-846-8270C	bis(2-Chloroethoxy)methane	<1.5	ug/L		1.48	4.9333	7/30/2001	JF
SW-846-8270C	bis(2-Chloroethyl)ether	<1.9	ug/L		1.9	6.3333	7/30/2001	JF
SW-846-8270C	bis(2-Chloroisopropyl)ether	<1.5	ug/L		1.54	5.1333	7/30/2001	JF
SW-846-8270C	bis(2-Ethylhexyl)phthalate	<4.6	ug/L		4.6	15.333	7/30/2001	JF
SW-846-8270C	Butylbenzylphthalate	<1.8	ug/L		1.8	6	7/30/2001	JF
SW-846-8270C	Chrysene	<1.6	ug/L		1.56	5.2	7/30/2001	JF
SW-846-8270C	Di-n-butylphthalate	<1.5	ug/L		1.52	5.0666	7/30/2001	JF
SW-846-8270C	Di-n-octylphthalate	<1.5	ug/L		1.52	5.0666	7/30/2001	JF
SW-846-8270C	Dibenz(a,j)acridine	<1.4	ug/L		1.44	4.8	7/30/2001	JF
SW-846-8270C	Dibenzo(a,h)anthracene	<1.5	ug/L		1.5	5	7/30/2001	JF
SW-846-8270C	Dibenzofuran	<1.3	ug/L		1.28	4.2666	7/30/2001	JF
SW-846-8270C	Diethylphthalate	<1.3	ug/L		1.26	4.2	7/30/2001	JF
SW-846-8270C	Dimethylphthalate	<1.5	ug/L		1.52	5.0666	7/30/2001	JF
SW-846-8270C	Diphenylamine	<2.6	ug/L		2.6	8.6666	7/30/2001	JF
SW-846-8270C	Ethyl methanesulfonate	<1.7	ug/L		1.7	5.6666	7/30/2001	JF
SW-846-8270C	Fluoranthene	<2.0	ug/L		1.98	6.6	7/30/2001	JF
SW-846-8270C	Fluorene	<1.2	ug/L		1.2	4	7/30/2001	JF
SW-846-8270C	Hexachlorobenzene	<1.6	ug/L		1.58	5.2666	7/30/2001	JF
SW-846-8270C	Hexachlorobutadiene	<7.6	ug/L		7.6	25.333	7/30/2001	JF
SW-846-8270C	Hexachlorocyclopentadiene	<5.4	ug/L		5.4	18	7/30/2001	JF
SW-846-8270C	Hexachloroethane	<5.2	ug/L		5.2	17.333	7/30/2001	JF
SW-846-8270C	Indeno(1,2,3-cd)pyrene	<1.4	ug/L		1.44	4.8	7/30/2001	JF
SW-846-8270C	Isophorone	<1.4	ug/L		1.42	4.7333	7/30/2001	JF
SW-846-8270C	Methyl methanesulfonate	<1.1	ug/L		1.06	3.5333	7/30/2001	JF
SW-846-8270C	n-Nitrosodi-n-butylamine	<1.5	ug/L		1.46	4.8666	7/30/2001	JF
SW-846-8270C	n-Nitrosodi-n-propylamine	<1.4	ug/L		1.4	4.6666	7/30/2001	JF
SW-846-8270C	n-Nitrosodimethylamine	<1.9	ug/L		1.88	6.2666	7/30/2001	JF
SW-846-8270C	n-Nitrosodiphenylamine	<2.6	ug/L		2.6	8.6666	7/30/2001	JF
SW-846-8270C	n-Nitrosopiperidine	<1.4	ug/L		1.4	4.6666	7/30/2001	JF
SW-846-8270C	Naphthalene	<1.4	ug/L		1.38	4.6	7/30/2001	JF
SW-846-8270C	Nitrobenzene	<1.4	ug/L		1.42	4.7333	7/30/2001	JF
SW-846-8270C	Nitrobenzene-d5 - Surrogate	51	% Rec		0	0	7/30/2001	JF
SW-846-8270C	p-Dimethylaminoazobenzene	<2.0	ug/L		2	6.6666	7/30/2001	JF
SW-846-8270C	Pentachlorobenzene	<2.2	ug/L		2.2	7.3333	7/30/2001	JF
SW-846-8270C	Pentachloronitrobenzene	<1.5	ug/L		1.54	5.1333	7/30/2001	JF
SW-846-8270C	Pentachlorophenol	<1.7	ug/L		1.72	5.7333	7/30/2001	JF
SW-846-8270C	Phenacetin	<1.8	ug/L		1.78	5.9333	7/30/2001	JF
SW-846-8270C	Phenanthrene	<1.4	ug/L		1.42	4.7333	7/30/2001	JF
SW-846-8270C	Phenol	<1.1	ug/L		1.1	3.6666	7/30/2001	JF
SW-846-8270C	Phenol-d5 - Surrogate	41	% Rec		0	0	7/30/2001	JF
SW-846-8270C	Pronamide	<1.7	ug/L		1.68	5.6	7/30/2001	JF
SW-846-8270C	Pyrene	<2.2	ug/L		2.2	7.3333	7/30/2001	JF
SW-846-8270C	Pyridine	<3.0	ug/L		3	10	7/30/2001	JF
SW-846-8270C	Terphenyl-d14 - Surrogate	59	% Rec		0	0	7/30/2001	JF

01REL010561 6/27/2001 DPZ-4

Robert E Lee & Associates, Inc
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 Project Number: 13551002
 Project Name: WEISENBERGER TIE & LUMBER

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 Fax: (920)336-9141
 Client ID: L14
 Chain: 85968
 Report Date: 8/23/2001

Method	Parameter Name	Result	Units	Flag	MDL	PQL	Anl Date	Analyst
Lab No.	Collect Date	Sample ID						
SW-846-8270C	Extraction Date	Complete			0	0	7/03/2001	JF
SW-846-8270C	1,2,4,5-Tetrachlorobenzene	<3.0	ug/L		3	10	7/30/2001	JF
SW-846-8270C	1,2,4-Trichlorobenzene	<3.6	ug/L		3.6	12	7/30/2001	JF
SW-846-8270C	1,2-Dichlorobenzene	<3.2	ug/L		3.2	10.666	7/30/2001	JF
SW-846-8270C	1,2-Diphenylhydrazine	<1.3	ug/L		1.34	4.4666	7/30/2001	JF
SW-846-8270C	1,3-Dichlorobenzene	<3.6	ug/L		3.6	12	7/30/2001	JF
SW-846-8270C	1,4-Dichlorobenzene	<3.4	ug/L		3.4	11.333	7/30/2001	JF
SW-846-8270C	1-Chloronaphthalene	<1.8	ug/L		1.84	6.1333	7/30/2001	JF
SW-846-8270C	1-Naphthylamine	<2.2	ug/L		2.2	7.3333	7/30/2001	JF
SW-846-8270C	2,3,4,6-Tetrachlorophenol	<1.4	ug/L		1.36	4.5333	7/30/2001	JF
SW-846-8270C	2,4,5-Trichlorophenol	<1.6	ug/L		1.64	5.4666	7/30/2001	JF
SW-846-8270C	2,4,6-Tribromophenol - Surrogate	71	% Rec		0	0	7/30/2001	JF
SW-846-8270C	2,4,6-Trichlorophenol	<1.6	ug/L		1.62	5.4	7/30/2001	JF
SW-846-8270C	2,4-Dichlorophenol	<1.5	ug/L		1.54	5.1333	7/30/2001	JF
SW-846-8270C	2,4-Dimethylphenol	<2.8	ug/L		2.8	9.3333	7/30/2001	JF
SW-846-8270C	2,4-Dinitrophenol	<1.8	ug/L		1.78	5.9333	7/30/2001	JF
SW-846-8270C	2,4-Dinitrotoluene	<1.6	ug/L		1.58	5.2666	7/30/2001	JF
SW-846-8270C	2,6-Dichlorophenol	<1.5	ug/L		1.52	5.0666	7/30/2001	JF
SW-846-8270C	2,6-Dinitrotoluene	<1.6	ug/L		1.6	5.3333	7/30/2001	JF
SW-846-8270C	2-Chloronaphthalene	<1.6	ug/L		1.56	5.2	7/30/2001	JF
SW-846-8270C	2-Chlorophenol	<1.2	ug/L		1.2	4	7/30/2001	JF
SW-846-8270C	2-Fluorobiphenyl - Surrogate	66	% Rec		0	0	7/30/2001	JF
SW-846-8270C	2-Fluorophenol - Surrogate	72	% Rec		0	0	7/30/2001	JF
SW-846-8270C	2-Methylnaphthalene	<1.6	ug/L		1.6	5.3333	7/30/2001	JF
SW-846-8270C	2-Methylphenol	<1.7	ug/L		1.7	5.6666	7/30/2001	JF
SW-846-8270C	2-Naphthylamine	<2.2	ug/L		2.2	7.3333	7/30/2001	JF
SW-846-8270C	2-Nitroaniline	<1.6	ug/L		1.64	5.4666	7/30/2001	JF
SW-846-8270C	2-Nitrophenol	<1.3	ug/L		1.28	4.2666	7/30/2001	JF
SW-846-8270C	2-Picoline	<3.0	ug/L		3	10	7/30/2001	JF
SW-846-8270C	3 & 4-Methylphenol	<1.8	ug/L		1.76	5.8666	7/30/2001	JF
SW-846-8270C	3,3'-Dichlorobenzidine	<0.84	ug/L		0.84	2.8	7/30/2001	JF
SW-846-8270C	3-Methylcholanthrene	<1.9	ug/L		1.94	6.4666	7/30/2001	JF
SW-846-8270C	3-Nitroaniline	<2.2	ug/L		2.2	7.3333	7/30/2001	JF
SW-846-8270C	4,6-Dinitro-2-methylphenol	<2.6	ug/L		2.6	8.6666	7/30/2001	JF
SW-846-8270C	4-Aminobiphenyl	<1.5	ug/L		1.5	5	7/30/2001	JF
SW-846-8270C	4-Bromophenyl phenyl ether	<1.2	ug/L		1.2	4	7/30/2001	JF
SW-846-8270C	4-Chloro-3-methylphenol	<1.7	ug/L		1.7	5.6666	7/30/2001	JF
SW-846-8270C	4-Chloroaniline	<1.3	ug/L		1.32	4.4	7/30/2001	JF
SW-846-8270C	4-Chlorophenyl phenyl ether	<1.2	ug/L		1.24	4.1333	7/30/2001	JF
SW-846-8270C	4-Nitroaniline	<1.3	ug/L		1.32	4.4	7/30/2001	JF
SW-846-8270C	4-Nitrophenol	<1.9	ug/L		1.86	6.2	7/30/2001	JF
SW-846-8270C	7,12-Dimethylbenz(a)anthracene	<2.0	ug/L		2	6.6666	7/30/2001	JF
SW-846-8270C	Acenaphthene	<1.1	ug/L		1.12	3.7333	7/30/2001	JF
SW-846-8270C	Acenaphthylene	<1.5	ug/L		1.46	4.8666	7/30/2001	JF
SW-846-8270C	Acetophenone	<1.6	ug/L		1.6	5.3333	7/30/2001	JF
SW-846-8270C	Aniline	<1.7	ug/L		1.74	5.8	7/30/2001	JF

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 Client ID: L14
 Chain: 85968
 Report Date: 8/23/2001

Method	Parameter Name	Result	Units	Flag	MDL	PQL	Anl Date	Analyst	
Lab No	Collect Date	Sample ID							
SW-846-8270C		Anthracene	<1.3		ug/L	1.3	4.3333	7/30/2001	JF
SW-846-8270C		Benzidine	<2.8		ug/L	2.8	9.3333	7/30/2001	JF
SW-846-8270C		Benzo(a)anthracene	<1.4		ug/L	1.42	4.7333	7/30/2001	JF
SW-846-8270C		Benzo(a)pyrene	<1.1		ug/L	1.1	3.6666	7/30/2001	JF
SW-846-8270C		Benzo(b)fluoranthene	<1.3		ug/L	1.28	4.2666	7/30/2001	JF
SW-846-8270C		Benzo(g,h,i)perylene	<1.3		ug/L	1.32	4.4	7/30/2001	JF
SW-846-8270C		Benzo(k)fluoranthene	<1.6		ug/L	1.58	5.2666	7/30/2001	JF
SW-846-8270C		Benzyl alcohol	<1.8		ug/L	1.8	6	7/30/2001	JF
SW-846-8270C		bis(2-Chloroethoxy)methane	<1.5		ug/L	1.48	4.9333	7/30/2001	JF
SW-846-8270C		bis(2-Chloroethyl)ether	<1.9		ug/L	1.9	6.3333	7/30/2001	JF
SW-846-8270C		bis(2-Chloroisopropyl)ether	<1.5		ug/L	1.54	5.1333	7/30/2001	JF
SW-846-8270C		bis(2-Ethylhexyl)phthalate	<4.6		ug/L	4.6	15.333	7/30/2001	JF
SW-846-8270C		Butylbenzylphthalate	<1.8		ug/L	1.8	6	7/30/2001	JF
SW-846-8270C		Chrysene	<1.6		ug/L	1.56	5.2	7/30/2001	JF
SW-846-8270C		Di-n-butylphthalate	<1.5		ug/L	1.52	5.0666	7/30/2001	JF
SW-846-8270C		Di-n-octylphthalate	<1.5		ug/L	1.52	5.0666	7/30/2001	JF
SW-846-8270C		Dibenz(a,j)acridine	<1.4		ug/L	1.44	4.8	7/30/2001	JF
SW-846-8270C		Dibenzo(a,h)anthracene	<1.5		ug/L	1.5	5	7/30/2001	JF
SW-846-8270C		Dibenzofuran	<1.3		ug/L	1.28	4.2666	7/30/2001	JF
SW-846-8270C		Diethylphthalate	<1.3		ug/L	1.26	4.2	7/30/2001	JF
SW-846-8270C		Dimethylphthalate	<1.5		ug/L	1.52	5.0666	7/30/2001	JF
SW-846-8270C		Diphenylamine	<2.6		ug/L	2.6	8.6666	7/30/2001	JF
SW-846-8270C		Ethyl methanesulfonate	<1.7		ug/L	1.7	5.6666	7/30/2001	JF
SW-846-8270C		Fluoranthene	<2.0		ug/L	1.98	6.6	7/30/2001	JF
SW-846-8270C		Fluorene	<1.2		ug/L	1.2	4	7/30/2001	JF
SW-846-8270C		Hexachlorobenzene	<1.6		ug/L	1.58	5.2666	7/30/2001	JF
SW-846-8270C		Hexachlorobutadiene	<7.6		ug/L	7.6	25.333	7/30/2001	JF
SW-846-8270C		Hexachlorocyclopentadiene	<5.4		ug/L	5.4	18	7/30/2001	JF
SW-846-8270C		Hexachloroethane	<5.2		ug/L	5.2	17.333	7/30/2001	JF
SW-846-8270C		Indeno(1,2,3-cd)pyrene	<1.4		ug/L	1.44	4.8	7/30/2001	JF
SW-846-8270C		Isophorone	<1.4		ug/L	1.42	4.7333	7/30/2001	JF
SW-846-8270C		Methyl methanesulfonate	<1.1		ug/L	1.06	3.5333	7/30/2001	JF
SW-846-8270C		n-Nitrosodi-n-butylamine	<1.5		ug/L	1.46	4.8666	7/30/2001	JF
SW-846-8270C		n-Nitrosodi-n-propylamine	<1.4		ug/L	1.4	4.6666	7/30/2001	JF
SW-846-8270C		n-Nitrosodimethylamine	<1.9		ug/L	1.88	6.2666	7/30/2001	JF
SW-846-8270C		n-Nitrosodiphenylamine	<2.6		ug/L	2.6	8.6666	7/30/2001	JF
SW-846-8270C		n-Nitrosopiperidine	<1.4		ug/L	1.4	4.6666	7/30/2001	JF
SW-846-8270C		Naphthalene	<1.4		ug/L	1.38	4.6	7/30/2001	JF
SW-846-8270C		Nitrobenzene	<1.4		ug/L	1.42	4.7333	7/30/2001	JF
SW-846-8270C		Nitrobenzene-d5 - Surrogate	59		% Rec	0	0	7/30/2001	JF
SW-846-8270C		p-Dimethylaminoazobenzene	<2.0		ug/L	2	6.6666	7/30/2001	JF
SW-846-8270C		Pentachlorobenzene	<2.2		ug/L	2.2	7.3333	7/30/2001	JF
SW-846-8270C		Pentachloronitrobenzene	<1.5		ug/L	1.54	5.1333	7/30/2001	JF
SW-846-8270C		Pentachlorophenol	5.2		ug/L	1.72	5.7333	7/30/2001	JF
SW-846-8270C		Phenacetin	<1.8		ug/L	1.78	5.9333	7/30/2001	JF
SW-846-8270C		Phenanthrene	<1.4		ug/L	1.42	4.7333	7/30/2001	JF

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Client ID: L14

Chain: 85968

Report Date: 8/23/2001

Method	Parameter Name	Result	Units	Flag	MDL	PQL	Anl.Date	Analyst
Lab No	Collect Date	Sample ID						
SW-846-8270C	Phenol	<1.1	ug/L		1.1	3.6666	7/30/2001	JF
SW-846-8270C	Phenol-d5 - Surrogate	39	% Rec		0	0	7/30/2001	JF
SW-846-8270C	Pronamide	<1.7	ug/L		1.68	5.6	7/30/2001	JF
SW-846-8270C	Pyrene	<2.2	ug/L		2.2	7.3333	7/30/2001	JF
SW-846-8270C	Pyridine	<3.0	ug/L		3	10	7/30/2001	JF
SW-846-8270C	Terphenyl-d14 - Surrogate	60	% Rec		0	0	7/30/2001	JF
01REL010562	6/27/2001	DPZ-5						
SW-846-8270C	Extraction Date	Complete			0	0	7/03/2001	JF
SW-846-8270C	1,2,4,5-Tetrachlorobenzene	<3.0	ug/L		3	10	7/30/2001	JF
SW-846-8270C	1,2,4-Trichlorobenzene	<3.6	ug/L		3.6	12	7/30/2001	JF
SW-846-8270C	1,2-Dichlorobenzene	<3.2	ug/L		3.2	10.666	7/30/2001	JF
SW-846-8270C	1,2-Diphenylhydrazine	<1.3	ug/L		1.34	4.4666	7/30/2001	JF
SW-846-8270C	1,3-Dichlorobenzene	<3.6	ug/L		3.6	12	7/30/2001	JF
SW-846-8270C	1,4-Dichlorobenzene	<3.4	ug/L		3.4	11.333	7/30/2001	JF
SW-846-8270C	1-Chloronaphthalene	<1.8	ug/L		1.84	6.1333	7/30/2001	JF
SW-846-8270C	1-Naphthylamine	<2.2	ug/L		2.2	7.3333	7/30/2001	JF
SW-846-8270C	2,3,4,6-Tetrachlorophenol	<1.4	ug/L		1.36	4.5333	7/30/2001	JF
SW-846-8270C	2,4,5-Trichlorophenol	<1.6	ug/L		1.64	5.4666	7/30/2001	JF
SW-846-8270C	2,4,6-Tribromophenol - Surrogate	57	% Rec		0	0	7/30/2001	JF
SW-846-8270C	2,4,6-Trichlorophenol	<1.6	ug/L		1.62	5.4	7/30/2001	JF
SW-846-8270C	2,4-Dichlorophenol	<1.5	ug/L		1.54	5.1333	7/30/2001	JF
SW-846-8270C	2,4-Dimethylphenol	<2.8	ug/L		2.8	9.3333	7/30/2001	JF
SW-846-8270C	2,4-Dinitrophenol	<1.8	ug/L		1.78	5.9333	7/30/2001	JF
SW-846-8270C	2,4-Dinitrotoluene	<1.6	ug/L		1.58	5.2666	7/30/2001	JF
SW-846-8270C	2,6-Dichlorophenol	<1.5	ug/L		1.52	5.0666	7/30/2001	JF
SW-846-8270C	2,6-Dinitrotoluene	<1.6	ug/L		1.6	5.3333	7/30/2001	JF
SW-846-8270C	2-Chloronaphthalene	<1.6	ug/L		1.56	5.2	7/30/2001	JF
SW-846-8270C	2-Chlorophenol	<1.2	ug/L		1.2	4	7/30/2001	JF
SW-846-8270C	2-Fluorobiphenyl - Surrogate	50	% Rec		0	0	7/30/2001	JF
SW-846-8270C	2-Fluorophenol - Surrogate	59	% Rec		0	0	7/30/2001	JF
SW-846-8270C	2-Methylnaphthalene	<1.6	ug/L		1.6	5.3333	7/30/2001	JF
SW-846-8270C	2-Methylphenol	<1.7	ug/L		1.7	5.6666	7/30/2001	JF
SW-846-8270C	2-Naphthylamine	<2.2	ug/L		2.2	7.3333	7/30/2001	JF
SW-846-8270C	2-Nitroaniline	<1.6	ug/L		1.64	5.4666	7/30/2001	JF
SW-846-8270C	2-Nitrophenol	<1.3	ug/L		1.28	4.2666	7/30/2001	JF
SW-846-8270C	2-Picoline	<3.0	ug/L		3	10	7/30/2001	JF
SW-846-8270C	3 & 4-Methylphenol	<1.8	ug/L		1.76	5.8666	7/30/2001	JF
SW-846-8270C	3,3'-Dichlorobenzidine	<0.84	ug/L		0.84	2.8	7/30/2001	JF
SW-846-8270C	3-Methylcholanthrene	<1.9	ug/L		1.94	6.4666	7/30/2001	JF
SW-846-8270C	3-Nitroaniline	<2.2	ug/L		2.2	7.3333	7/30/2001	JF
SW-846-8270C	4,6-Dinitro-2-methylphenol	<2.6	ug/L		2.6	8.6666	7/30/2001	JF
SW-846-8270C	4-Aminobiphenyl	<1.5	ug/L		1.5	5	7/30/2001	JF
SW-846-8270C	4-Bromophenyl phenyl ether	<1.2	ug/L		1.2	4	7/30/2001	JF
SW-846-8270C	4-Chloro-3-methylphenol	<1.7	ug/L		1.7	5.6666	7/30/2001	JF
SW-846-8270C	4-Chloroaniline	<1.3	ug/L		1.32	4.4	7/30/2001	JF
SW-846-8270C	4-Chlorophenyl phenyl ether	<1.2	ug/L		1.24	4.1333	7/30/2001	JF

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Method	Parameter Name	Result	Units	Flag	MDL	PQL	Anl Date	Analyst
Lab No.	Collect Date	Sample ID						
SW-846-8270C	4-Nitroaniline	<1.3	ug/L		1.32	4.4	7/30/2001	JF
SW-846-8270C	4-Nitrophenol	<1.9	ug/L		1.86	6.2	7/30/2001	JF
SW-846-8270C	7,12-Dimethylbenz(a)anthracene	<2.0	ug/L		2	6.6666	7/30/2001	JF
SW-846-8270C	Acenaphthene	<1.1	ug/L		1.12	3.7333	7/30/2001	JF
SW-846-8270C	Acenaphthylene	<1.5	ug/L		1.46	4.8666	7/30/2001	JF
SW-846-8270C	Acetophenone	<1.6	ug/L		1.6	5.3333	7/30/2001	JF
SW-846-8270C	Aniline	<1.7	ug/L		1.74	5.8	7/30/2001	JF
SW-846-8270C	Anthracene	<1.3	ug/L		1.3	4.3333	7/30/2001	JF
SW-846-8270C	Benzidine	<2.8	ug/L		2.8	9.3333	7/30/2001	JF
SW-846-8270C	Benzo(a)anthracene	<1.4	ug/L		1.42	4.7333	7/30/2001	JF
SW-846-8270C	Benzo(a)pyrene	<1.1	ug/L		1.1	3.6666	7/30/2001	JF
SW-846-8270C	Benzo(b)fluoranthene	<1.3	ug/L		1.28	4.2666	7/30/2001	JF
SW-846-8270C	Benzo(g,h,i)perylene	<1.3	ug/L		1.32	4.4	7/30/2001	JF
SW-846-8270C	Benzo(k)fluoranthene	<1.6	ug/L		1.58	5.2666	7/30/2001	JF
SW-846-8270C	Benzyl alcohol	<1.8	ug/L		1.8	6	7/30/2001	JF
SW-846-8270C	bis(2-Chloroethoxy)methane	<1.5	ug/L		1.48	4.9333	7/30/2001	JF
SW-846-8270C	bis(2-Chloroethyl)ether	<1.9	ug/L		1.9	6.3333	7/30/2001	JF
SW-846-8270C	bis(2-Chloroisopropyl)ether	<1.5	ug/L		1.54	5.1333	7/30/2001	JF
SW-846-8270C	bis(2-Ethylhexyl)phthalate	<4.6	ug/L		4.6	15.333	7/30/2001	JF
SW-846-8270C	Butylbenzylphthalate	<1.8	ug/L		1.8	6	7/30/2001	JF
SW-846-8270C	Chrysene	<1.6	ug/L		1.56	5.2	7/30/2001	JF
SW-846-8270C	Di-n-butylphthalate	<1.5	ug/L		1.52	5.0666	7/30/2001	JF
SW-846-8270C	Di-n-octylphthalate	<1.5	ug/L		1.52	5.0666	7/30/2001	JF
SW-846-8270C	Dibenz(a,j)acridine	<1.4	ug/L		1.44	4.8	7/30/2001	JF
SW-846-8270C	Dibenzo(a,h)anthracene	<1.5	ug/L		1.5	5	7/30/2001	JF
SW-846-8270C	Dibenzofuran	<1.3	ug/L		1.28	4.2666	7/30/2001	JF
SW-846-8270C	Diethylphthalate	<1.3	ug/L		1.26	4.2	7/30/2001	JF
SW-846-8270C	Dimethylphthalate	<1.5	ug/L		1.52	5.0666	7/30/2001	JF
SW-846-8270C	Diphenylamine	<2.6	ug/L		2.6	8.6666	7/30/2001	JF
SW-846-8270C	Ethyl methanesulfonate	<1.7	ug/L		1.7	5.6666	7/30/2001	JF
SW-846-8270C	Fluoranthene	<2.0	ug/L		1.98	6.6	7/30/2001	JF
SW-846-8270C	Fluorene	<1.2	ug/L		1.2	4	7/30/2001	JF
SW-846-8270C	Hexachlorobenzene	<1.6	ug/L		1.58	5.2666	7/30/2001	JF
SW-846-8270C	Hexachlorobutadiene	<7.6	ug/L		7.6	25.333	7/30/2001	JF
SW-846-8270C	Hexachlorocyclopentadiene	<5.4	ug/L		5.4	18	7/30/2001	JF
SW-846-8270C	Hexachloroethane	<5.2	ug/L		5.2	17.333	7/30/2001	JF
SW-846-8270C	Indeno(1,2,3-cd)pyrene	<1.4	ug/L		1.44	4.8	7/30/2001	JF
SW-846-8270C	Isophorone	<1.4	ug/L		1.42	4.7333	7/30/2001	JF
SW-846-8270C	Methyl methanesulfonate	<1.1	ug/L		1.06	3.5333	7/30/2001	JF
SW-846-8270C	n-Nitrosodi-n-butylamine	<1.5	ug/L		1.46	4.8666	7/30/2001	JF
SW-846-8270C	n-Nitrosodi-n-propylamine	<1.4	ug/L		1.4	4.6666	7/30/2001	JF
SW-846-8270C	n-Nitrosodimethylamine	<1.9	ug/L		1.88	6.2666	7/30/2001	JF
SW-846-8270C	n-Nitrosodiphenylamine	<2.6	ug/L		2.6	8.6666	7/30/2001	JF
SW-846-8270C	n-Nitrosopiperidine	<1.4	ug/L		1.4	4.6666	7/30/2001	JF
SW-846-8270C	Naphthalene	<1.4	ug/L		1.38	4.6	7/30/2001	JF
SW-846-8270C	Nitrobenzene	<1.4	ug/L		1.42	4.7333	7/30/2001	JF

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 Fax: (920)336-9141
 Client ID: L14
 Chain: 85968
 Report Date: 8/23/2001

Method	Parameter Name	Result	Units	Flag	MDL	PQL	Anl Date	Analyst
Lab No.	Collect Date	Sample ID						
SW-846-8270C		Nitrobenzene-d5 - Surrogate	44	% Rec	0	0	7/30/2001	JF
SW-846-8270C		p-Dimethylaminoazobenzene	<2.0	ug/L	2	6.6666	7/30/2001	JF
SW-846-8270C		Pentachlorobenzene	<2.2	ug/L	2.2	7.3333	7/30/2001	JF
SW-846-8270C		Pentachloronitrobenzene	<1.5	ug/L	1.54	5.1333	7/30/2001	JF
SW-846-8270C		Pentachlorophenol	<1.7	ug/L	1.72	5.7333	7/30/2001	JF
SW-846-8270C		Phenacetin	<1.8	ug/L	1.78	5.9333	7/30/2001	JF
SW-846-8270C		Phenanthrene	<1.4	ug/L	1.42	4.7333	7/30/2001	JF
SW-846-8270C		Phenol	<1.1	ug/L	1.1	3.6666	7/30/2001	JF
SW-846-8270C		Phenol-d5 - Surrogate	34	% Rec	0	0	7/30/2001	JF
SW-846-8270C		Pronamide	<1.7	ug/L	1.68	5.6	7/30/2001	JF
SW-846-8270C		Pyrene	<2.2	ug/L	2.2	7.3333	7/30/2001	JF
SW-846-8270C		Pyridine	<3.0	ug/L	3	10	7/30/2001	JF
SW-846-8270C		Terphenyl-d14 - Surrogate	48	% Rec	0	0	7/30/2001	JF

01REL010563 6/27/2001 DPZ-6

EPA 1613	Dioxin Analysis	See Attached			0	0	7/25/2001	PJK
SW-846-8270C	Extraction Date	Complete			0	0	7/03/2001	JF
SW-846-8270C	1,2,4,5-Tetrachlorobenzene	<3.0	ug/L		3	10	7/29/2001	JF
SW-846-8270C	1,2,4-Trichlorobenzene	<3.6	ug/L		3.6	12	7/29/2001	JF
SW-846-8270C	1,2-Dichlorobenzene	<3.2	ug/L		3.2	10.666	7/29/2001	JF
SW-846-8270C	1,2-Diphenylhydrazine	<1.3	ug/L		1.34	4.4666	7/29/2001	JF
SW-846-8270C	1,3-Dichlorobenzene	<3.6	ug/L		3.6	12	7/29/2001	JF
SW-846-8270C	1,4-Dichlorobenzene	<3.4	ug/L		3.4	11.333	7/29/2001	JF
SW-846-8270C	1-Chloronaphthalene	<1.8	ug/L		1.84	6.1333	7/29/2001	JF
SW-846-8270C	1-Naphthylamine	<2.2	ug/L		2.2	7.3333	7/29/2001	JF
SW-846-8270C	2,3,4,6-Tetrachlorophenol	<1.4	ug/L		1.36	4.5333	7/29/2001	JF
SW-846-8270C	2,4,5-Trichlorophenol	<1.6	ug/L		1.64	5.4666	7/29/2001	JF
SW-846-8270C	2,4,6-Tribromophenol - Surrogate	62	% Rec		0	0	7/29/2001	JF
SW-846-8270C	2,4,6-Trichlorophenol	<1.6	ug/L		1.62	5.4	7/29/2001	JF
SW-846-8270C	2,4-Dichlorophenol	<1.5	ug/L		1.54	5.1333	7/29/2001	JF
SW-846-8270C	2,4-Dimethylphenol	<2.8	ug/L		2.8	9.3333	7/29/2001	JF
SW-846-8270C	2,4-Dinitrophenol	<1.8	ug/L		1.78	5.9333	7/29/2001	JF
SW-846-8270C	2,4-Dinitrotoluene	<1.6	ug/L		1.58	5.2666	7/29/2001	JF
SW-846-8270C	2,6-Dichlorophenol	<1.5	ug/L		1.52	5.0666	7/29/2001	JF
SW-846-8270C	2,6-Dinitrotoluene	<1.6	ug/L		1.6	5.3333	7/29/2001	JF
SW-846-8270C	2-Chloronaphthalene	<1.6	ug/L		1.56	5.2	7/29/2001	JF
SW-846-8270C	2-Chlorophenol	<1.2	ug/L		1.2	4	7/29/2001	JF
SW-846-8270C	2-Fluorobiphenyl - Surrogate	57	% Rec		0	0	7/29/2001	JF
SW-846-8270C	2-Fluorophenol - Surrogate	44	% Rec		0	0	7/29/2001	JF
SW-846-8270C	2-Methylnaphthalene	<1.6	ug/L		1.6	5.3333	7/29/2001	JF
SW-846-8270C	2-Methylphenol	<1.7	ug/L		1.7	5.6666	7/29/2001	JF
SW-846-8270C	2-Naphthylamine	<2.2	ug/L		2.2	7.3333	7/29/2001	JF
SW-846-8270C	2-Nitroaniline	<1.6	ug/L		1.64	5.4666	7/29/2001	JF
SW-846-8270C	2-Nitrophenol	<1.3	ug/L		1.28	4.2666	7/29/2001	JF
SW-846-8270C	2-Picoline	<3.0	ug/L		3	10	7/29/2001	JF
SW-846-8270C	3 & 4-Methylphenol	<1.8	ug/L		1.76	5.8666	7/29/2001	JF
SW-846-8270C	3,3'-Dichlorobenzidine	<0.84	ug/L		0.84	2.8	7/29/2001	JF

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 Client ID: L14
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 Report Date: 8/23/2001

Method	Parameter Name	Result	Units	Flag	MDL	PQL	Anl Date	Analyst
Lab No	Collect Date	Sample ID						
SW-846-8270C	3-Methylcholanthrene	<1.9	ug/L		1.94	6.4666	7/29/2001	JF
SW-846-8270C	3-Nitroaniline	<2.2	ug/L		2.2	7.3333	7/29/2001	JF
SW-846-8270C	4,6-Dinitro-2-methylphenol	<2.6	ug/L		2.6	8.6666	7/29/2001	JF
SW-846-8270C	4-Aminobiphenyl	<1.5	ug/L		1.5	5	7/29/2001	JF
SW-846-8270C	4-Bromophenyl phenyl ether	<1.2	ug/L		1.2	4	7/29/2001	JF
SW-846-8270C	4-Chloro-3-methylphenol	<1.7	ug/L		1.7	5.6666	7/29/2001	JF
SW-846-8270C	4-Chloroaniline	<1.3	ug/L		1.32	4.4	7/29/2001	JF
SW-846-8270C	4-Chlorophenyl phenyl ether	<1.2	ug/L		1.24	4.1333	7/29/2001	JF
SW-846-8270C	4-Nitroaniline	<1.3	ug/L		1.32	4.4	7/29/2001	JF
SW-846-8270C	4-Nitrophenol	<1.9	ug/L		1.86	6.2	7/29/2001	JF
SW-846-8270C	7,12-Dimethylbenz(a)anthracene	<2.0	ug/L		2	6.6666	7/29/2001	JF
SW-846-8270C	Acenaphthene	<1.1	ug/L		1.12	3.7333	7/29/2001	JF
SW-846-8270C	Acenaphthylene	<1.5	ug/L		1.46	4.8666	7/29/2001	JF
SW-846-8270C	Acetophenone	<1.6	ug/L		1.6	5.3333	7/29/2001	JF
SW-846-8270C	Aniline	<1.7	ug/L		1.74	5.8	7/29/2001	JF
SW-846-8270C	Anthracene	<1.3	ug/L		1.3	4.3333	7/29/2001	JF
SW-846-8270C	Benzidine	<2.8	ug/L		2.8	9.3333	7/29/2001	JF
SW-846-8270C	Benzo(a)anthracene	<1.4	ug/L		1.42	4.7333	7/29/2001	JF
SW-846-8270C	Benzo(a)pyrene	<1.1	ug/L		1.1	3.6666	7/29/2001	JF
SW-846-8270C	Benzo(b)fluoranthene	<1.3	ug/L		1.28	4.2666	7/29/2001	JF
SW-846-8270C	Benzo(g,h,i)perylene	<1.3	ug/L		1.32	4.4	7/29/2001	JF
SW-846-8270C	Benzo(k)fluoranthene	<1.6	ug/L		1.58	5.2666	7/29/2001	JF
SW-846-8270C	Benzyl alcohol	<1.8	ug/L		1.8	6	7/29/2001	JF
SW-846-8270C	bis(2-Chloroethoxy)methane	<1.5	ug/L		1.48	4.9333	7/29/2001	JF
SW-846-8270C	bis(2-Chloroethyl)ether	<1.9	ug/L		1.9	6.3333	7/29/2001	JF
SW-846-8270C	bis(2-Chloroisopropyl)ether	<1.5	ug/L		1.54	5.1333	7/29/2001	JF
SW-846-8270C	bis(2-Ethylhexyl)phthalate	<4.6	ug/L		4.6	15.333	7/29/2001	JF
SW-846-8270C	Butylbenzylphthalate	<1.8	ug/L		1.8	6	7/29/2001	JF
SW-846-8270C	Chrysene	<1.6	ug/L		1.56	5.2	7/29/2001	JF
SW-846-8270C	Di-n-butylphthalate	<1.5	ug/L		1.52	5.0666	7/29/2001	JF
SW-846-8270C	Di-n-octylphthalate	<1.5	ug/L		1.52	5.0666	7/29/2001	JF
SW-846-8270C	Dibenz(a,j)acridine	<1.4	ug/L		1.44	4.8	7/29/2001	JF
SW-846-8270C	Dibenzo(a,h)anthracene	<1.5	ug/L		1.5	5	7/29/2001	JF
SW-846-8270C	Dibenzofuran	<1.3	ug/L		1.28	4.2666	7/29/2001	JF
SW-846-8270C	Diethylphthalate	<1.3	ug/L		1.26	4.2	7/29/2001	JF
SW-846-8270C	Dimethylphthalate	<1.5	ug/L		1.52	5.0666	7/29/2001	JF
SW-846-8270C	Diphenylamine	<2.6	ug/L		2.6	8.6666	7/29/2001	JF
SW-846-8270C	Ethyl methanesulfonate	<1.7	ug/L		1.7	5.6666	7/29/2001	JF
SW-846-8270C	Fluoranthene	<2.0	ug/L		1.98	6.6	7/29/2001	JF
SW-846-8270C	Fluorene	<1.2	ug/L		1.2	4	7/29/2001	JF
SW-846-8270C	Hexachlorobenzene	<1.6	ug/L		1.58	5.2666	7/29/2001	JF
SW-846-8270C	Hexachlorobutadiene	<7.6	ug/L		7.6	25.333	7/29/2001	JF
SW-846-8270C	Hexachlorocyclopentadiene	<5.4	ug/L		5.4	18	7/29/2001	JF
SW-846-8270C	Hexachloroethane	<5.2	ug/L		5.2	17.333	7/29/2001	JF
SW-846-8270C	Indeno(1,2,3-cd)pyrene	<1.4	ug/L		1.44	4.8	7/29/2001	JF
SW-846-8270C	Isophorone	<1.4	ug/L		1.42	4.7333	7/29/2001	JF

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Method	Parameter Name	Result	Units	Flag	MDL	PQL	Anl Date	Analyst
Lab No.	Collect Date	Sample ID						
SW-846-8270C	Methyl methanesulfonate	<1.1	ug/L		1.06	3.5333	7/29/2001	JF
SW-846-8270C	n-Nitrosodi-n-butylamine	<1.5	ug/L		1.46	4.8666	7/29/2001	JF
SW-846-8270C	n-Nitrosodi-n-propylamine	<1.4	ug/L		1.4	4.6666	7/29/2001	JF
SW-846-8270C	n-Nitrosodimethylamine	<1.9	ug/L		1.88	6.2666	7/29/2001	JF
SW-846-8270C	n-Nitrosodiphenylamine	<2.6	ug/L		2.6	8.6666	7/29/2001	JF
SW-846-8270C	n-Nitrosopiperidine	<1.4	ug/L		1.4	4.6666	7/29/2001	JF
SW-846-8270C	Naphthalene	<1.4	ug/L		1.38	4.6	7/29/2001	JF
SW-846-8270C	Nitrobenzene	<1.4	ug/L		1.42	4.7333	7/29/2001	JF
SW-846-8270C	Nitrobenzene-d5 - Surrogate	51	% Rec		0	0	7/29/2001	JF
SW-846-8270C	p-Dimethylaminoazobenzene	<2.0	ug/L		2	6.6666	7/29/2001	JF
SW-846-8270C	Pentachlorobenzene	<2.2	ug/L		2.2	7.3333	7/29/2001	JF
SW-846-8270C	Pentachloronitrobenzene	<1.5	ug/L		1.54	5.1333	7/29/2001	JF
SW-846-8270C	Pentachlorophenol	<1.7	ug/L		1.72	5.7333	7/29/2001	JF
SW-846-8270C	Phenacetin	<1.8	ug/L		1.78	5.9333	7/29/2001	JF
SW-846-8270C	Phenanthrene	<1.4	ug/L		1.42	4.7333	7/29/2001	JF
SW-846-8270C	Phenol	<1.1	ug/L		1.1	3.6666	7/29/2001	JF
SW-846-8270C	Phenol-d5 - Surrogate	39	% Rec		0	0	7/29/2001	JF
SW-846-8270C	Pronamide	<1.7	ug/L		1.68	5.6	7/29/2001	JF
SW-846-8270C	Pyrene	<2.2	ug/L		2.2	7.3333	7/29/2001	JF
SW-846-8270C	Pyridine	<3.0	ug/L		3	10	7/29/2001	JF
SW-846-8270C	Terphenyl-d14 - Surrogate	50	% Rec		0	0	7/29/2001	JF
<u>01REL010564 6/27/2001 DUP 1</u>								
SW-846-8021B	1,2,4-Trimethylbenzene	81	ug/L		0.23	0.7666	7/10/2001	CRW
SW-846-8021B	1,3,5-Trimethylbenzene	17	ug/L		0.21	0.7	7/10/2001	CRW
SW-846-8021B	Benzene	3.4	ug/L		0.21	0.7	7/10/2001	CRW
SW-846-8021B	Ethylbenzene	12	ug/L		0.23	0.7666	7/10/2001	CRW
SW-846-8021B	Fluorobenzene-Surrogate	97	% Rec		0	0	7/10/2001	CRW
SW-846-8021B	Methyl-tertiary-butyl ether	<0.091	ug/L		0.091	0.3033	7/10/2001	CRW
SW-846-8021B	Toluene	2.2	ug/L		0.22	0.7333	7/10/2001	CRW
SW-846-8021B	Xylenes-Total	73	ug/L		0.44	1.4666	7/10/2001	CRW
<u>01REL010565 6/27/2001 EQUIP B1</u>								
SW-846-8021B	1,2,4-Trimethylbenzene	<0.23	ug/L		0.23	0.7666	7/10/2001	CRW
SW-846-8021B	1,3,5-Trimethylbenzene	<0.21	ug/L		0.21	0.7	7/10/2001	CRW
SW-846-8021B	Benzene	<0.21	ug/L		0.21	0.7	7/10/2001	CRW
SW-846-8021B	Ethylbenzene	<0.23	ug/L		0.23	0.7666	7/10/2001	CRW
SW-846-8021B	Fluorobenzene-Surrogate	101	% Rec		0	0	7/10/2001	CRW
SW-846-8021B	Methyl-tertiary-butyl ether	<0.091	ug/L		0.091	0.3033	7/10/2001	CRW
SW-846-8021B	Toluene	<0.22	ug/L		0.22	0.7333	7/10/2001	CRW
SW-846-8021B	Xylenes-Total	<0.44	ug/L		0.44	1.4666	7/10/2001	CRW
<u>01REL010566 6/27/2001 TRIP B</u>								
SW-846-8021B	1,2,4-Trimethylbenzene	<0.23	ug/L		0.23	0.7666	7/10/2001	CRW
SW-846-8021B	1,3,5-Trimethylbenzene	<0.21	ug/L		0.21	0.7	7/10/2001	CRW
SW-846-8021B	Benzene	<0.21	ug/L		0.21	0.7	7/10/2001	CRW
SW-846-8021B	Ethylbenzene	<0.23	ug/L		0.23	0.7666	7/10/2001	CRW
SW-846-8021B	Fluorobenzene-Surrogate	100	% Rec		0	0	7/10/2001	CRW

Robert E. Lee & Associates, Inc.
 Wisconsin Certification Number: 405043870
 Certificate of Analysis Report

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Method	Parameter Name	Result	Units	Flag	MDL	PQL	Anl>Date	Analyst
Lab No.	Collect Date	Sample ID						
SW-846-8021B	Methyl-tertiary-butyl ether	<0.091	ug/L		0.091	0.3033	7/10/2001	CRW
SW-846-8021B	Toluene	<0.22	ug/L		0.22	0.7333	7/10/2001	CRW
SW-846-8021B	Xylenes-Total	<0.44	ug/L		0.44	1.4666	7/10/2001	CRW
01REL010567	6/27/2001	BK 859						
EPA-515.1	Chlorinated Acids Pesticides Analysis	See Attached			0	0	8/07/2001	PJK

Robert E. Lee & Associates. Inc.

Quality Control Report - Description of Flags

Flag	Section	Description
13	L	The reported result is less than the practical quantitation limit (PQL).
21	L	Reported result is above calibration curve.

WEISENBERGER TIE & LUMBER
PROJECT # 3551-002
MARCH 2001
GROUNDWATER SAMPLING

WELLS: ~~MW-3+~~ DMW-3+ ~~DMW-4+~~ ~~DMW-5+~~ ~~EQUIP B1+~~
~~DMW-6A+~~ ~~DMW-10+~~ ~~DPZ-1+~~ ~~DUP-1+~~ ~~TRIP B+~~

PARAMETERS: PVOCs - EPA METHOD 8020

SAMPLE AMOUNT: 2-40ML VIALS - HCL

WELLS: ~~MW-3+~~ ~~MW-5+~~ ~~MW-6+~~ ~~MW-+~~ ~~MW-10+~~
~~DMW-1+~~ ~~DMW-2+~~ DMW-3+ ~~DMW-4+~~ ~~DMW-5+~~
~~DMW-6A+~~ ~~DMW-7+~~ ~~DMW-10+~~ ~~DMW-13+~~ ~~DPZ-1+~~
~~DPZ-1A+~~ ~~DPZ-2+~~ ~~DPZ-3+~~ ~~DPZ-4+~~ ~~DPZ-5+~~ ~~DPZ-6+~~

PARAMETERS: BNA - EPA METHOD 8270

SAMPLE AMOUNT: 1 LITER AMBER - UNPRESERVED

WELLS: ~~MW-3-~~ ~~MW-5-2~~ ~~MW-6-~~ ~~MW-7-~~ ~~MW-10-~~ ~~DMW-1-~~
~~DMW-2-~~ DMW-3- ~~DMW-4-~~ ~~DMW-5-~~ ~~DMW-6A-~~ ~~DMW-7-~~
~~DMW-8-~~ ~~DPZ-1-~~ ~~DPZ-3-~~ ~~DPZ-6-~~

PARAMETERS: BNA - DIOXIN/FURANS - EPA METHOD 8290

SAMPLE AMOUNT: 1 LITER AMBER - UNPRESERVED

PRIVATE WELL - KRAUTKRAMER - BARN WELL

SAMPLE ID: BK859-

PARAMETERS: PENTACHLOROPHENOL - EPA METHOD 515

SAMPLE AMOUNT: 1 LITER AMBER - HCL

We assume DMW-3 is dry as we received no bottle.



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Milwaukee Office 262.569.8893 FAX 262.569.7995

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CHAIN OF CUSTODY RECORD

coc # 85968 (10f3)

Client: <u>Weisenberger Tie + Lumber</u>		Project Number: <u>1355100</u>		Analyses Required: (Note special detection limits or methods)		Report to:			
Project Name: <u>1</u>				No. Of Containers Preservation Type (see key below) <u>see enclosed sheet</u>		Company:			
PO #: <u>1</u>		BID #: <u>2</u>				Address:			
Environmental Program: <input type="checkbox"/> LUST <input type="checkbox"/> SDWA <input type="checkbox"/> WPDES <input type="checkbox"/> RCRA <input type="checkbox"/> OTHER		Requested Turnaround Time: <input checked="" type="checkbox"/> Normal (10-15 DAYS) <input type="checkbox"/> Rush				Check Delivery Method: <input checked="" type="checkbox"/> In Person <input type="checkbox"/> Mail <input type="checkbox"/> Common Courier <input type="checkbox"/> Courier Service <input type="checkbox"/> Other		Telephone:	
Date Needed: _____		Rushes accepted only w/prior notification				Invoice To:		Company:	
Sampler: <u>Craig W...</u>		Sample Type (Matrix): DW = Drinking Water GW = Groundwater WW = Wastewater Soil, Oil, Sludge, Air, Other:				Address:			
Sample Name	Date	Time	Comp	Lab	Pres	REL Sample No.	Remarks:		
MW-3	6/27/01	9:00		X	N	6W	4 U, H X	10543 Puoc, Bna, Diox	
MW-5		3:15	A				2 U	10544 Bna, Diox	
MW-6			A				2 U	10545 Bna, Diox	
MW-7			A				2 U	10546 Bna, Diox	
MW-10			A				2 U	10547 Bna, Diox	
DMW-1			A				2 U	10548 Bna, Diox	
DMW-2			A				4 2 U	10549 Bna, Diox	
DMW-4			A				4 U, H	10550 Puoc, Bna, Diox	
DMW-5			A				4 U, H	10551 Puoc, Bna, Diox	
DMW-6A			A				2 U, H	10552 Bna, Diox, Puoc	
DMW-7			A				2 U	10553 Bna, Diox	
DMW-8			A				1 U	10554 Diox	
Relinquished By: <u>Craig W...</u>		Date: <u>6/27/01</u>	Time: <u>6:00</u>	Received By: _____		Date: _____	Time: _____	Laboratory Receiving Notes Temperature of Contents _____ °C Custody Seal Intact _____ Sample Condition _____ Sample pH _____	
Received by Lab: <u>Merrill Amire</u>									
1) _____		A/P							
2) _____		A/P							
3) _____		A/P							

WISCONSIN DNR CERTIFICATION NUMBER 405043870

Preservation Key

N = Nitric Acid O = Sodium Hydroxide
 H = Hydrochloric Acid U = Unpreserved
 M = Methanol S = Sulfuric Acid



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 Milwaukee Office 262.569.8893 FAX 262.569.7995

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CHAIN OF CUSTODY RECORD

COC # **85968 (2 of 3)**

Client: Wessenberg, Tel and Jmbr		Project Name: 1		Project Number: Y3551002		Analyses Required: (Note special detection limits or methods)		Report to:		
PO #: 1		BID #:		Environmental Program: <input type="checkbox"/> LUST <input type="checkbox"/> SDWA <input type="checkbox"/> WPDES <input type="checkbox"/> RCRA <input type="checkbox"/> OTHER				Company:		Address:
Requested Turnaround Time <input checked="" type="checkbox"/> Normal (10-15 DAYS) <input type="checkbox"/> Rush		Check Delivery Method <input checked="" type="checkbox"/> In Person <input type="checkbox"/> Mail <input type="checkbox"/> Common Courier <input type="checkbox"/> Courier Service <input type="checkbox"/> Other		No. of Containers		Preservation Type (see key below)		Telephone:		
Date Needed: _____ <small>Rushes accepted only w/prior notification</small>		Sampler: Craig W...						Sample Type (Matrix) DW = Drinking Water GW = Groundwater WW = Wastewater Soil, Oil, Sludge, Air, Other:		Fax:
Sample Name	Date	Time	Comp	Lab	Pres	Mat	No. of Containers	Preservation Type	REL Sample No.	Remarks
DMW-10	6/27/01	9:00	A	X	N	GW	3	U	10557	Puoc, Bna
DMW-13	6/27/01	3:15	A				1	U	10556	Bna
DPZ-1			A				4	OH	10555	Puoc, Bna, Diox
DPZ-1A			A				1	U	10558	Bna
DPZ-2			A				1	U	10559	Bna
DPZ-3			A				2	U	10560	Bna, Diox
DPZ-4			A				1	U	10561	Bna
DPZ-5			A				1	U	10562	Bna
DPZ-6			A				2	U	10563	Bna, Diox
Dup 1			A				4	H	10564	Puoc
Range B 1			A				2	H	10565	Puoc
Trig B			A				2	H	10566	Puoc

Relinquished By: Craig W...	Date: 6/27/01	Time: 6:00	Received By: _____	Date: _____	Time: _____
1) _____			A/P		
2) _____			A/P		
3) _____			A/P		
Received by Lab: Herry Ferrero	Date: 6/27/01				

WISCONSIN DNR CERTIFICATION NUMBER 405043870

Preservation Key
 N = Nitric Acid O = Sodium Hydroxide
 H = Hydrochloric Acid U = Unpreserved
 M = Methanol S = Sulfuric Acid



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CHAIN OF CUSTODY RECORD

COC #

259108 (30f3)

Client: <u>Wasserman Tie + lumber</u>				Analyses Required: (Note special detection limits or methods)				Report to:			
Project Name: <u>1</u>		Project Number: <u>1355 1002</u>						Company:			
PO #: <u>1</u>		BID #: _____						Address:			
Environmental Program: <input type="checkbox"/> LUST <input type="checkbox"/> SDWA <input type="checkbox"/> WPDES <input type="checkbox"/> RCRA <input type="checkbox"/> OTHER _____								Telephone:			
Requested Turnaround Time: <input checked="" type="checkbox"/> Normal (10-15 DAYS) <input type="checkbox"/> Rush		Check Delivery Method: <input checked="" type="checkbox"/> In Person <input type="checkbox"/> Mail <input type="checkbox"/> Common Courier <input type="checkbox"/> Courier Service <input type="checkbox"/> Other _____		No. of Containers Preservation Type (see key below) <u>see enclosed sheet</u>		Fax:		Invoice To:			
Date Needed: _____ <small>Rushes accepted only w/prior notification</small>		Sampler: <u>[Signature]</u>				Company:		Address:			
Sample Name		Date				Time		Telephone:		Fax:	
								REL Sample No.		Remarks:	
<u>BR 859</u>		<u>6/27/01</u>		<u>11:30</u>		<u>10577</u>		<u>515</u>			

Relinquished By: <u>[Signature]</u>		Date: <u>6/27/01</u>		Time: <u>6:00</u>		Received By: _____		Date: _____		Time: _____	
1) _____		_____		_____		_____		_____		A/P	
2) _____		_____		_____		_____		_____		A/P	
3) _____		_____		_____		_____		_____		A/P	
Received by Lab: <u>[Signature]</u>				<u>6/27/01</u>							

Laboratory Receiving Notes

Temperature of Contents _____ °C

Custody Seal Intact _____

Sample Condition _____

Sample pH _____

A = AM P = PM

Preservation Key

N = Nitric Acid O = Sodium Hydroxide
H = Hydrochloric Acid U = Unpreserved
M = Methanol S = Sulfuric Acid

**LABORATORY REPORT
PACE ANALYTICAL LABORATORY
DIOXIN/FURAN ANALYSIS**

- **MONITORING WELLS**

DETERMINATION OF PCDD/PCDF LEVELS

Prepared for:
Robert E. Lee & Associates, Inc.
Attn: Paul Knuth
2825 South Webster Avenue
Box 2100
Green Bay, WI 54306

Project: Chemical Analysis

Client Purchase Order Number: NA

REPORT OF LABORATORY ANALYSIS

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REPORT OF: CHEMICAL ANALYSES

PROJECT: PCDD/PCDF ANALYSES

DATE: July 30, 2001

ISSUED TO: Robert E. Lee & Associates, Inc.
Attn: Mr. Paul Knuth
2825 South Webster Avenue
Box 2100
Green Bay, WI 54306

REPORT NO:01-1046199

INTRODUCTION

This report presents the results from the analyses performed on fifteen samples which were submitted by a representative of Robert E. Lee & Associates, Inc. The samples were analyzed for the presence or absence of polychlorinated dibenzo-p-dioxins (PCDDs) and dibenzofurans (PCDFs) using a modified version of USEPA Method 8290 as described below.

SAMPLE IDENTIFICATION

<u>Client ID</u>	<u>Sample Type</u>	<u>Date Received</u>	<u>Pace ID</u>
01-10543	Water	06/29/01	102831328
01-10544	Water	06/29/01	102831336
01-10545	Water	06/29/01	102831344
01-10546	Water	06/29/01	102831351
01-10547	Water	06/29/01	102831369
01-10548	Water	06/29/01	102831377
01-10549	Water	06/29/01	102831385
01-10550	Water	06/29/01	102831393
01-10551	Water	06/29/01	102831401
01-10552	Water	06/29/01	102831419
01-10553	Water	06/29/01	102831427
01-10554	Water	06/29/01	102831435
01-10557	Water	06/29/01	102831443
01-10560	Water	06/29/01	102831450
01-10563	Water	06/29/01	102831468

REPORT OF LABORATORY ANALYSIS

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PROJECT: PCDD/PCDF ANALYSES

DATE: July 30, 2001

PAGE: 2

REPORT NO 01-1046199

METHODOLOGY

Sample Extraction

Each sample was spiked with $^{13}\text{C}_{12}$ -labeled PCDD/PCDF internal standards (Table 1) and extracted with methylene chloride in a separatory funnel. The extract was quantitatively transferred to a Kuderna-Danish concentrator, concentrated, and solvent exchanged to hexane. The hexane extract was then spiked with 2,3,7,8-TCDD- $^{37}\text{Cl}_4$ enrichment efficiency standard (Table 1) and processed through the analyte enrichment procedures described below.

PCDD/PCDF Analyte Enrichment

The extraction procedure often removes a variety of compounds, in addition to the PCDDs and PCDFs, from the sample matrix. Some of these compounds can directly interfere with the analyses while others can overload the capillary column causing degradation in chromatographic resolution or sensitivity. The analyte enrichment steps described below are used to remove interferences from the extracts.

Each extract was diluted to 100 mL with hexane, transferred to a separatory funnel, and washed with 1N sodium hydroxide, concentrated sulfuric acid, and aqueous sodium chloride (5% w/v) as needed. The hexane extract was quantitatively transferred to a liquid chromatography column containing alternating layers of silica gel, 40% concentrated sulfuric acid on silica gel, and 33% 1 N sodium hydroxide on silica gel. The column was eluted with 90 mL of hexane and the entire eluate was collected and concentrated, under ambient conditions, to a volume of 1 mL and spiked with the $^{37}\text{Cl}_4$ -TCDD cleanup standard (Table 1).

Each extract was then fractionated on a liquid chromatography column containing 4 g of activated alumina. The column was eluted with 20 mL of hexane followed by 15 mL of 60% methylene chloride/hexane. The 60% methylene chloride/hexane fraction was collected, concentrated, spiked with recovery standards (1,2,3,4-TCDD- $^{13}\text{C}_{12}$ and 1,2,3,7,8,9-HxCDD- $^{13}\text{C}_{12}$) and taken to a final volume of 20 μL .

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REPORT OF: CHEMICAL ANALYSES

PROJECT: PCDD/PCDF ANALYSES

DATE: July 30, 2001

PAGE: 3

REPORT NO: 01-1046199

PCDD/PCDF Analyses

Each sample extract was analyzed for the presence of PCDDs and PCDFs using combined capillary column gas chromatography/high resolution mass spectrometry (HRGC/HRMS). The instrumentation consisted of a Hewlett Packard Model 6890 gas chromatograph interfaced to a Micromass Ultima high-resolution mass spectrometer. The capillary column was interfaced directly into the ion source of the mass spectrometer, thus providing the highest possible sensitivity while minimizing degradation of the chromatographic resolution.

The mass spectrometer was operated in the electron impact ionization mode at a mass resolution of 10,000-11,000 ($M/\Delta M$, 10 percent valley definition). This resolution is sufficient to resolve most interferences, such as PCBs, thus providing the highest level of confidence that the detected levels of PCDD/PCDF were not false positives resulting from interferences. Typical operating parameters for the HRGC/HRMS analyses are summarized in Table 2.

The data were acquired by selected-ion-recording (SIR) using groups of ion masses similar to those described in USEPA Method 8290. The five groups corresponded to the tetrachlorinated through octachlorinated congener classes. Each group contained two ion masses for the PCDDs, two ion masses for the PCDFs, the corresponding ion masses from the two isotopically labeled internal standards, and the ion mass characteristic of the polychlorinated diphenylether (PCDE) which, if present, could cause false responses in the dibenzofuran channels.

Each group of ion masses also contained a lock mass which was used by the data system to automatically correct the mass focus of the instrument. The data system determined the centroid of the lock mass during each data acquisition cycle and corrected the mass focus of the analyte and internal standard ion masses to assure that the centers of the mass peaks were being monitored.

The criteria used to judge positive responses for a PCDD/PCDF isomer included:

- * Simultaneous response at both ion masses of the PCDD or PCDF
- * Signal-to-noise ratio equal to or greater than 2.5:1.0 for both ion masses
- * Chlorine isotope ratio within 15% of the theoretical value
- * Chromatographic retention time within +/- 2 seconds of the expected retention time
- * Chromatographic retention times within elution windows determined from analyses of standard mixtures
- * Absence of simultaneous response in the PCDF and PCDE ion traces

A list of the exact ion masses monitored for the determination of PCDD/PCDF isomers and the PCDE interferences is presented in Table 3. Also included are the theoretical chlorine isotope ratios for the ten congener classes. **REPORT OF LABORATORY ANALYSIS**

PROJECT: PCDD/PCDF ANALYSES

DATE: July 30, 2001

PAGE: 4

REPORT NO: 01-1046199

PCDD/PCDF Quantification and Calculations

The PCDD/PCDF isomers were quantified by comparison of their responses to the responses of the labeled internal standards. Relative response factors were calculated from analyses of standard mixtures containing representatives of each of the PCDD/PCDF congener classes at five concentration levels, and each of the internal standards at one concentration level, as shown in Table 4. The PCDD/PCDF response factors were calculated by comparing the sum of the responses from the two ion masses monitored for each chlorine congener class to the sum of the responses from the two ion masses of the corresponding isotopically labeled internal standard. The formula for the response factor calculation is:

$$R_f = \frac{A_n \times Q_{is}}{A_{is} \times Q_n}$$

where:

- Rf = Response factor
- A_n = Sum of integrated areas for native isomer
- Q_{is} = Quantity of labeled internal standard
- A_{is} = Sum of integrated areas for labeled internal standard
- Q_n = Quantity of native isomer

The levels of PCDD/PCDF in each sample were quantified using the following equation:

$$C = \frac{A_n \times Q_{is}}{A_{is} \times W \times R_f}$$

where:

- C = Concentration of target isomer or congener class
- A_n = Sum of integrated areas for the target isomer or congener class
- Q_{is} = Quantity of labeled internal standard added to the sample
- A_{is} = Sum of integrated areas for the labeled internal standard
- W = Sample amount
- Rf = Response factor

REPORT OF LABORATORY ANALYSIS

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REPORT OF: CHEMICAL ANALYSES

PROJECT: PCDD/PCDF ANALYSES

DATE: July 30, 2001

PAGE: 5

REPORT NO: 01-1046199

PCDD/PCDF Quantification and Calculations (Cont.)

Each pair of ion mass peaks in the selected-ion-current chromatograms was evaluated manually to determine if it met the criteria for a PCDD or PCDF isomer. Areas of all peaks exhibiting correct ion ratios, having retention times within the correct windows, and having areas corresponding to concentrations in the range covered by the initial calibration were then summed for calculations of total congener concentrations.

A limit of detection (LOD) based on producing a signal that is 2.5 times the noise level, was calculated for each undetected 2,3,7,8-substituted isomer of any tetra through octa chlorinated congener class. The noise heights used to calculate the detection limits were measured at the retention time of the specific isomer. The formula used for calculating the LOD is:

$$\text{LOD} = \frac{H_n \times Q_{is} \times 2.5}{H_{is} \times W \times R_f}$$

where:

LOD = Single isomer limit of detection
H_n = Sum of noise heights at native isomer retention time
Q_{is} = Quantity of labeled internal standard
H_{is} = Sum of peak heights for labeled internal standard
W = Sample amount
R_f = Response factor

The recovery of the 2,3,7,8-TCDD-³⁷Cl₄ enrichment efficiency standard and each ¹³C₁₂-labeled internal standard, relative to either 1,2,3,4-TCDD-¹³C₁₂ or 1,2,3,7,8,9-HxCDD-¹³C₁₂, was calculated using the following equation:

$$\%R = \frac{A_{is} \times Q_{rs} \times 100\%}{R_{fr} \times A_{rs} \times Q_{is}}$$

where:

%R = Percent recovery of labeled internal standard
A_{is} = Sum of integrated areas of labeled internal standard
Q_{rs} = Quantity of recovery standard
A_{rs} = Sum of integrated areas of recovery standard
R_{fr} = Response factor of the specific labeled internal standard relative to the recovery standard
Q_{is} = Quantity of the labeled internal standard congener added to the sample

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REPORT OF: CHEMICAL ANALYSES

PROJECT: PCDD/PCDF ANALYSES

DATE: July 30, 2001

PAGE: 6

REPORT NO: 01-1046199

Quality Control for PCDD/PCDF Analyses

The performance of the sample processing steps and the instrumentation are monitored on a routine basis. The procedures and criteria are summarized below.

One method blank and one laboratory spike sample are typically prepared with each ten samples of any given matrix. Recoveries of the native PCDD/PCDF analytes in the laboratory spike samples generally range from 70 to 130%. Recoveries of selected analytes outside this range do not invalidate the data but provide information, which is used by the laboratory to monitor recovery trends and to assure optimization of the method.

Internal standards are spiked into each sample prior to extraction in order to monitor the level of recovery, which is achieved for each individual sample. Acceptable recoveries range from 40 to 135 percent for the internal standards unless a deviation is due to variation in instrument response as a result of analytical interferences.

The resolution of the mass spectrometer is verified prior to each analysis to be 10,000 or greater. Hardcopies of the reference peaks are printed at the beginning and end of each analysis day. The resolving power of the DB-5MS chromatographic column is checked daily by analyzing a standard solution containing 2,3,7,8-TCDD and the adjacent TCDD isomers. The DB-225 column resolution is checked daily by analyzing a standard solution containing 2,3,7,8-TCDF and the adjacent TCDF isomers. Acceptable performance is achieved when 2,3,7,8-TCDD or 2,3,7,8-TCDF is resolved from the adjacent isomers by a valley of 25% or less. The group times for the selected-ion-monitoring data acquisitions are also checked daily by analyzing the column performance mix which has been modified to contain the first and last eluting isomers of each congener class. In this way one is assured of collecting data representative of the total PCDD/PCDF content and that the 2,3,7,8-substituted isomers are suitably resolved.

Initial calibrations are generated by analyzing standard solutions (see Table 4) containing target native and labeled PCDD/PCDF compounds. Response factors are calculated and averaged for each compound. These averages are used for quantification and for comparison to the daily continuing calibration. The relative standard deviation for each native compound must be 20% or less (30% or less for the labeled compounds) as specified in Method 8290. A continuing calibration standard is analyzed at the beginning and end of each 12-hour shift on days when initial calibrations are not performed. The initial calibration is considered to be valid when the response factors from the continuing calibration analysis fall to within the ranges specified in Method 8290.

REPORT OF LABORATORY ANALYSIS

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REPORT OF: CHEMICAL ANALYSES

PROJECT: PCDD/PCDF ANALYSES

DATE: July 30, 2001

PAGE: 7

REPORT NO: 01-1046199

RESULTS

The results from the analyses are presented in the following:

- Appendix A - Documentation
- Appendix B - PCDD/PCDF Analysis Results

DISCUSSION

The recoveries of the isotopically-labeled PCDD/PCDF internal standards in the sample extracts generally ranged from 59-132% and indicate a level of efficiency through the extraction and enrichment steps that is considered typical for this matrix. With the exception of four internal standards, which were affected by an interference, the labeled standard recoveries for the samples were within the Method 8290 target ranges. Also, since the quantifications of the native 2,3,7,8-substituted isomers were based on isotope dilution, the data were automatically corrected for variation in recovery and accurate values were obtained.

Some of the samples were found to contain compounds which interfere with the determination of co-eluting PCDD and PCDF isomers. Any affected 2,3,7,8-substituted isomers are flagged "E" or "I" on the data summary sheets.

A laboratory method blank was prepared and analyzed with the sample batch as part of our routine quality control procedures. The results, found at the beginning of Appendix B, show the blank to contain a trace level of OCDD. The blank sample level was below the calibration range of the method. Two of the samples contained this isomer at a level similar to that seen in the blank and are flagged "B" on the data summary sheet. In general, levels less than ten times the background are not considered statistically different from the background.

Laboratory spike samples were prepared with the sample batch by extracting laboratory water that had been fortified with native standard materials. Recoveries of the native compounds in the spiked samples ranged from 86-105% with relative percent differences of 1.0-17.0%. This indicates high degrees of accuracy and precision for these determinations.

REPORT OF LABORATORY ANALYSIS

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PROJECT: PCDD/PCDF ANALYSES

DATE: July 30, 2001

PAGE: 8

REPORT NO: 01-1046199

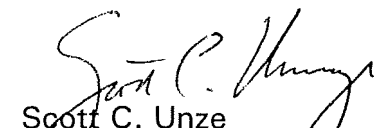
REMARKS

The sample extracts will be retained for a period of 30 days from the date of this report and then discarded unless other arrangements are made. The raw mass spectral data will be archived on magnetic tape for a period of not less than one year. Questions regarding the data contained in this report may be directed to the authors at the numbers provided below.

Pace Analytical Services, Inc.



Charles V. Sueper, Technical Director
High Resolution Mass Spectrometry
(612) 607-6387



Scott C. Unze
Project Manager, Dioxins
(612) 607-6383

REPORT OF LABORATORY ANALYSIS

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TABLE 1. Spike Levels of PCDD/PCDF Standards

Internal Standards	Spike Level (ng)
2,3,7,8-TCDF- ¹³ C ₁₂	2.0
2,3,7,8-TCDD- ¹³ C ₁₂	2.0
1,2,3,7,8-PeCDF- ¹³ C ₁₂	2.0
2,3,4,7,8-PeCDF- ¹³ C ₁₂	2.0
1,2,3,7,8-PeCDD- ¹³ C ₁₂	2.0
1,2,3,4,7,8-HxCDF- ¹³ C ₁₂	2.0
1,2,3,6,7,8-HxCDF- ¹³ C ₁₂	2.0
1,2,3,7,8,9-HxCDF- ¹³ C ₁₂	2.0
2,3,4,6,7,8-HxCDF- ¹³ C ₁₂	2.0
1,2,3,4,7,8-HxCDD- ¹³ C ₁₂	2.0
1,2,3,6,7,8-HxCDD- ¹³ C ₁₂	2.0
1,2,3,4,6,7,8-HpCDF- ¹³ C ₁₂	2.0
1,2,3,4,7,8,9-HpCDF- ¹³ C ₁₂	2.0
1,2,3,4,6,7,8-HpCDD- ¹³ C ₁₂	2.0
OCDD- ¹³ C ₁₂	4.0
<u>Recovery Standards</u>	
1,2,3,4-TCDD- ¹³ C ₁₂	2.0
1,2,3,7,8,9-HxCDD- ¹³ C ₁₂	2.0
<u>Enrichment Efficiency Standard</u>	
2,3,7,8-TCDD- ³⁷ Cl ₄	0.2

REPORT OF LABORATORY ANALYSIS

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**TABLE 2. High Resolution PCDD/PCDF Analyses
HRGC/HRMS Operating Parameters**

Mass Resolution	10,000-11,000 (M/ Δ M, 10% valley)
Electron Energy	32 electron volts
Accelerating Voltage	8,000 volts
Source Temperature	275°C
Preamplifier Gain	10 ⁻⁶ amp/volt
Multiplier Gain	~ 10 ⁵
Chromatographic Column	60 M DB-5MS
Transfer Line Temperature	260°C
Injection Mode	Splitless
Carrier Gas	Helium
Carrier Flow Velocity	~ 30 cm/sec

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**TABLE 3. Exact Ion Masses Monitored
for the Determination of PCDDs, PCDFs, and PCDEs**

Ratio Compound	Accurate Mass		Theoretical
	Mass 1	Mass 2	Mass 1/Mass 2
Tetra-CDDs	319.8965	321.8936	0.77
Tetra-CDFs	303.9016	305.8987	0.77
Hexa-CDEs	375.8364		
Penta-CDDs	355.8546	357.8517	1.54
Penta-CDFs	339.8597	341.8567	1.54
Hepta-CDEs	409.7974		
Hexa-CDDs	389.8156	391.8127	1.23
Hexa-CDFs	373.8207	375.8178	1.23
Octa-CDEs	445.7555		
Hepta-CDDs	423.7766	425.7737	1.03
Hepta-CDFs	407.7817	409.7788	1.03
Nona-CDEs	479.7165		
Octa-CDD	457.7377	459.7347	0.88
Octa-CDF	441.7428	443.7398	0.88
Deca-CDE	513.6775		

CDDs = Chlorinated Dibenzo-p-dioxins
CDFs = Chlorinated Dibenzofurans
CDEs = Chlorinated Diphenylethers

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TABLE 4. High Resolution Calibration Solutions

Native CDDs/CDFs	Concentration (pg/uL)				
	CS1	CS2	CS3	CS4	CS5
2,3,7,8-TCDD	0.5	2	10	40	200
2,3,7,8 TCDF	0.5	2	10	40	200
1,2,3,7,8-PeCDD	2.5	10	50	200	1000
1,2,3,7,8-PeCDF	2.5	10	50	200	1000
2,3,4,7,8-PeCDF	2.5	10	50	200	1000
1,2,3,4,7,8-HxCDD	2.5	10	50	200	1000
1,2,3,6,7,8-HxCDD	2.5	10	50	200	1000
1,2,3,7,8,9-HxCDD	2.5	10	50	200	1000
1,2,3,4,7,8-HxCDF	2.5	10	50	200	1000
1,2,3,6,7,8-HxCDF	2.5	10	50	200	1000
1,2,3,7,8,9-HxCDF	2.5	10	50	200	1000
2,3,4,6,7,8-HxCDF	2.5	10	50	200	1000
1,2,3,4,6,7,8-HpCDD	2.5	10	50	200	1000
1,2,3,4,6,7,8-HpCDF	2.5	10	50	200	1000
1,2,3,4,7,8,9-HpCDF	2.5	10	50	200	1000
OCDD	5.0	20	100	400	2000
OCDF	5.0	20	100	400	2000
Internal Standards					
2,3,7,8-TCDD- ¹³ C ₁₂	100	100	100	100	100
2,3,7,8-TCDF- ¹³ C ₁₂	100	100	100	100	100
1,2,3,7,8-PeCDD- ¹³ C ₁₂	100	100	100	100	100
1,2,3,7,8-PeCDF- ¹³ C ₁₂	100	100	100	100	100
2,3,4,7,8-PeCDF- ¹³ C ₁₂	100	100	100	100	100
1,2,3,4,7,8-HxCDD- ¹³ C ₁₂	100	100	100	100	100
1,2,3,6,7,8-HxCDD- ¹³ C ₁₂	100	100	100	100	100
1,2,3,4,7,8-HxCDF- ¹³ C ₁₂	100	100	100	100	100
1,2,3,6,7,8-HxCDF- ¹³ C ₁₂	100	100	100	100	100
1,2,3,7,8,9-HxCDF- ¹³ C ₁₂	100	100	100	100	100
2,3,4,6,7,8-HxCDF- ¹³ C ₁₂	100	100	100	100	100
1,2,3,4,6,7,8-HpCDD- ¹³ C ₁₂	100	100	100	100	100
1,2,3,4,6,7,8-HpCDF- ¹³ C ₁₂	100	100	100	100	100
1,2,3,4,7,8,9-HpCDF- ¹³ C ₁₂	100	100	100	100	100
OCDD- ¹³ C ₁₂	200	200	200	200	200
Recovery Standards					
1,2,3,4-TCDD- ¹³ C ₁₂	100	100	100	100	100
1,2,3,7,8,9-HxCDD- ¹³ C ₁₂	100	100	100	100	100
Enrichment Efficiency Standard					
2,3,7,8-TCDD- ³⁷ C ₁₄	0.5	2	10	40	200

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TABLE 5. 2,3,7,8-TCDD Equivalency Factors (TEFs) for the Polychlorinated Dibenzo-p-dioxins and Dibenzofurans

Number	Compound(s)	TEF
1	2,3,7,8-TCDD	1.00
2	1,2,3,7,8-PeCDD	0.50
3	1,2,3,6,7,8-HxCDD	0.1
4	1,2,3,7,8,9-HxCDD	0.1
5	1,2,3,4,7,8-HxCDD	0.1
6	1,2,3,4,6,7,8-HpCDD	0.01
7	OCDD	0.001
8	* Total - TCDD	0.0
9	* Total - PeCDD	0.0
10	* Total - HxCDD	0.0
11	* Total - HpCDD	0.0
12	2,3,7,8-TCDF	0.10
13	1,2,3,7,8-PeCDF	0.05
14	2,3,4,7,8-PeCDF	0.5
15	1,2,3,6,7,8-HxCDF	0.1
16	1,2,3,7,8,9-HxCDF	0.1
17	1,2,3,4,7,8-HxCDF	0.1
18	2,3,4,6,7,8-HxCDF	0.1
19	1,2,3,4,6,7,8-HpCDF	0.01
20	1,2,3,4,7,8,9-HpCDF	0.01
21	OCDF	0.001
22	* Total - TCDF	0.0
23	* Total - PeCDF	0.0
24	* Total - HxCDF	0.0
25	* Total - HpCDF	0.0

*Excluding the 2,3,7,8-substituted congeners.

Reference: 1989 ITEFs

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Appendix A

REPORT OF LABORATORY ANALYSIS

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Appendix B

REPORT OF LABORATORY ANALYSIS

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Method 8290 Blank Analysis Results

Client - ROBERT E LEE

Lab Sample ID	BLANK-1153	Matrix	WATER
Filename	U10724A_10	Dilution	NA
Total Amount Extracted	1022.4 mL	Extracted	07/10/2001
ICAL Date	07/24/2001	Analyzed	07/24/2001 20:21
CCal Filename(s)	U10724A_06 & U10724A_21	Injected By	CVS

Native Isomers	Conc ng/L	EMPC ng/L	LRL ng/L		Internal Standards	ng's Added	Percent Recovery
2,3,7,8-TCDF	ND	----	0.0074	A	2,3,7,8-TCDF-13C	2.00	91
Total TCDF	ND	----	0.0020		2,3,7,8-TCDD-13C	2.00	86
					1,2,3,7,8-PeCDF-13C	2.00	93
2,3,7,8-TCDD	ND	----	0.0120	A	2,3,4,7,8-PeCDF-13C	2.00	92
Total TCDD	ND	----	0.0020		1,2,3,7,8-PeCDD-13C	2.00	91
					1,2,3,4,7,8-HxCDF-13C	2.00	93
1,2,3,7,8-PeCDF	ND	----	0.0098		1,2,3,6,7,8-HxCDF-13C	2.00	84
2,3,4,7,8-PeCDF	ND	----	0.0098		2,3,4,6,7,8-HxCDF-13C	2.00	99
Total PeCDF	ND	----	0.0098		1,2,3,7,8,9-HxCDF-13C	2.00	90
					1,2,3,4,7,8-HxCDD-13C	2.00	104
1,2,3,7,8-PeCDD	ND	----	0.0110	A	1,2,3,6,7,8-HxCDD-13C	2.00	96
Total PeCDD	ND	----	0.0098		1,2,3,4,6,7,8-HpCDF-13C	2.00	92
					1,2,3,4,7,8,9-HpCDF-13C	2.00	89
1,2,3,4,7,8-HxCDF	ND	----	0.0120	A	1,2,3,4,6,7,8-HpCDD-13C	2.00	93
1,2,3,6,7,8-HxCDF	ND	----	0.0098		OCDD-13C	4.00	80
2,3,4,6,7,8-HxCDF	ND	----	0.0098				
1,2,3,7,8,9-HxCDF	ND	----	0.0140	A	1,2,3,4-TCDD-13C	2.00	NA
Total HxCDF	ND	----	0.0098		1,2,3,7,8,9-HxCDD-13C	2.00	NA
1,2,3,4,7,8-HxCDD	ND	----	0.0180	A	2,3,7,8-TCDD-37Cl4	0.20	79
1,2,3,6,7,8-HxCDD	ND	----	0.0190	A			
1,2,3,7,8,9-HxCDD	ND	----	0.0220	A			
Total HxCDD	ND	----	0.0098				
1,2,3,4,6,7,8-HpCDF	----	0.013	0.0100	IA	Total 2,3,7,8-TCDD		
1,2,3,4,7,8,9-HpCDF	ND	----	0.0150	A	Equivalence: 0.000046 ng/L		
Total HpCDF	ND	----	0.0098		(Using ITE Factors)		
1,2,3,4,6,7,8-HpCDD	ND	----	0.0220	A			
Total HpCDD	ND	----	0.0098				
OCDF	ND	----	0.0230	A			
OCDD	0.046	----	0.0220	JA			

Conc = Concentration (Totals include 2,3,7,8-substituted isomers).
EMPC = Estimated Maximum Possible Concentration
LRL = Lower Reporting Limit
J = Concentration detected is below the calibration range
P = Recovery outside of target range
A = Detection Limit based on signal-to-noise measurement

I = Interference
E = PCDE Interference
ND = Not Detected
NA = Not Applicable
NC = Not Calculated
* = See Discussion

Report No.....01-1046199

REPORT OF LABORATORY ANALYSIS

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Method 8290 Analysis Results

Client - ROBERT E LEE

Client's Sample ID	01-10543		
Lab Sample ID	102831328		
Filename	U10721B_08		
Injected By	BAL		
Total Amount Extracted	1055.3 mL	Matrix	WATER
% Moisture	NA	Dilution	NA
Dry Weight Extracted	NA	Collected	06/27/2001
ICAL Date	07/21/2001	Received	06/29/2001
CCal Filename(s)	U10721B_01 & U10721B_16	Extracted	07/10/2001
Method Blank ID	BLANK-1153	Analyzed	07/22/2001 02:10

Native Isomers	Conc ng/L	EMPC ng/L	LRL ng/L		Internal Standards	ng's Added	Percent Recovery
2,3,7,8-TCDF	ND	----	0.0030	A	2,3,7,8-TCDF-13C	2.00	88
Total TCDF	ND	----	0.0019		2,3,7,8-TCDD-13C	2.00	84
					1,2,3,7,8-PeCDF-13C	2.00	105
2,3,7,8-TCDD	ND	----	0.0047	A	2,3,4,7,8-PeCDF-13C	2.00	114
Total TCDD	ND	----	0.0019		1,2,3,7,8-PeCDD-13C	2.00	103
					1,2,3,4,7,8-HxCDF-13C	2.00	113
1,2,3,7,8-PeCDF	ND	----	0.0095		1,2,3,6,7,8-HxCDF-13C	2.00	125
2,3,4,7,8-PeCDF	ND	----	0.0095		2,3,4,6,7,8-HxCDF-13C	2.00	117
Total PeCDF	ND	----	0.0095		1,2,3,7,8,9-HxCDF-13C	2.00	107
					1,2,3,4,7,8-HxCDD-13C	2.00	113
1,2,3,7,8-PeCDD	ND	----	0.0095		1,2,3,6,7,8-HxCDD-13C	2.00	113
Total PeCDD	ND	----	0.0095		1,2,3,4,6,7,8-HpCDF-13C	2.00	87
					1,2,3,4,7,8,9-HpCDF-13C	2.00	73
1,2,3,4,7,8-HxCDF	ND	----	0.0095		1,2,3,4,6,7,8-HpCDD-13C	2.00	75
1,2,3,6,7,8-HxCDF	ND	----	0.0095		OCDD-13C	4.00	61
2,3,4,6,7,8-HxCDF	ND	----	0.0095				
1,2,3,7,8,9-HxCDF	ND	----	0.0095		1,2,3,4-TCDD-13C	2.00	NA
Total HxCDF	0.043	----	0.0095	J	1,2,3,7,8,9-HxCDD-13C	2.00	NA
1,2,3,4,7,8-HxCDD	ND	----	0.0095		2,3,7,8-TCDD-37Cl4	0.20	86
1,2,3,6,7,8-HxCDD	----	0.012	0.0095	I			
1,2,3,7,8,9-HxCDD	ND	----	0.0095				
Total HxCDD	0.035	----	0.0095	J			
1,2,3,4,6,7,8-HpCDF	----	0.140	0.0095	E	Total 2,3,7,8-TCDD		
1,2,3,4,7,8,9-HpCDF	ND	----	0.0140	A	Equivalence: 0.0044 ng/L		
Total HpCDF	ND	----	0.0095		(Using ITE Factors)		
1,2,3,4,6,7,8-HpCDD	0.220	----	0.0140	A			
Total HpCDD	0.430	----	0.0095				
OCDF	0.170	----	0.0240	A			
OCDD	2.000	----	0.0280	A			

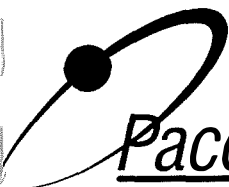
Conc = Concentration (Totals include 2,3,7,8-substituted isomers)
EMPC = Estimated Maximum Possible Concentration
A = Detection Limit based on signal-to-noise measurement
J = Concentration detected is below the calibration range
B = Less than 10 times higher than method blank level
P = Recovery outside of target range
Nn = Value obtained from additional analysis

LRL = Lower Reporting Limit
I = Interference
E = PCDE Interference
S = Saturated signal
ND = Not Detected
NA = Not Applicable
NC = Not Calculated
* = See Discussion

Report No.....01-1046199

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Method 8290 Analysis Results

Client - ROBERT E LEE

Client's Sample ID	01-10544		
Lab Sample ID	102831336		
Filename	U10721B_09		
Injected By	BAL		
Total Amount Extracted	1053.5 mL	Matrix	WATER
% Moisture	NA	Dilution	NA
Dry Weight Extracted	NA	Collected	06/27/2001
ICAL Date	07/21/2001	Received	06/29/2001
CCal Filename(s)	U10721B_01 & U10721B_16	Extracted	07/10/2001
Method Blank ID	BLANK-1153	Analyzed	07/22/2001 03:00

Native Isomers	Conc ng/L	EMPC ng/L	LRL ng/L		Internal Standards	ng's Added	Percent Recovery
2,3,7,8-TCDF	ND	----	0.0046	A	2,3,7,8-TCDF-13C	2.00	83
Total TCDF	ND	----	0.0019		2,3,7,8-TCDD-13C	2.00	77
					1,2,3,7,8-PeCDF-13C	2.00	97
2,3,7,8-TCDD	ND	----	0.0061	A	2,3,4,7,8-PeCDF-13C	2.00	107
Total TCDD	ND	----	0.0019		1,2,3,7,8-PeCDD-13C	2.00	97
					1,2,3,4,7,8-HxCDF-13C	2.00	117
1,2,3,7,8-PeCDF	ND	----	0.0095		1,2,3,6,7,8-HxCDF-13C	2.00	123
2,3,4,7,8-PeCDF	ND	----	0.0095		2,3,4,6,7,8-HxCDF-13C	2.00	128
Total PeCDF	ND	----	0.0095		1,2,3,7,8,9-HxCDF-13C	2.00	109
					1,2,3,4,7,8-HxCDD-13C	2.00	145 IP
1,2,3,7,8-PeCDD	ND	----	0.0095		1,2,3,6,7,8-HxCDD-13C	2.00	130
Total PeCDD	ND	----	0.0095		1,2,3,4,6,7,8-HpCDF-13C	2.00	93
					1,2,3,4,7,8,9-HpCDF-13C	2.00	77
1,2,3,4,7,8-HxCDF	ND	----	0.0095		1,2,3,4,6,7,8-HpCDD-13C	2.00	83
1,2,3,6,7,8-HxCDF	ND	----	0.0095		OCDD-13C	4.00	70
2,3,4,6,7,8-HxCDF	ND	----	0.0095				
1,2,3,7,8,9-HxCDF	ND	----	0.0095		1,2,3,4-TCDD-13C	2.00	NA
Total HxCDF	ND	----	0.0095		1,2,3,7,8,9-HxCDD-13C	2.00	NA
1,2,3,4,7,8-HxCDD	ND	----	0.0095		2,3,7,8-TCDD-37Cl4	0.20	74
1,2,3,6,7,8-HxCDD	ND	----	0.0095				
1,2,3,7,8,9-HxCDD	ND	----	0.0095				
Total HxCDD	ND	----	0.0095				
1,2,3,4,6,7,8-HpCDF	ND	----	0.0130	A	Total 2,3,7,8-TCDD		
1,2,3,4,7,8,9-HpCDF	ND	----	0.0110	A	Equivalence: 0.00013 ng/L		
Total HpCDF	ND	----	0.0095		(Using ITE Factors)		
1,2,3,4,6,7,8-HpCDD	ND	----	0.0160	A			
Total HpCDD	ND	----	0.0095				
OCDF	ND	----	0.0390	A			
OCDD	0.13	----	0.0550	BA			

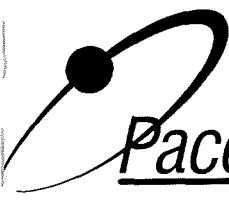
Conc = Concentration (Totals include 2,3,7,8-substituted isomers)
 EMPC = Estimated Maximum Possible Concentration
 A = Detection Limit based on signal-to-noise measurement
 J = Concentration detected is below the calibration range
 B = Less than 10 times higher than method blank level
 P = Recovery outside of target range
 Nn = Value obtained from additional analysis

LRL = Lower Reporting Limit
 I = Interference
 E = PCDE Interference
 S = Saturated signal
 ND = Not Detected
 NA = Not Applicable
 NC = Not Calculated
 * = See Discussion

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Method 8290 Analysis Results

Client - ROBERT E LEE

Client's Sample ID	01-10545			
Lab Sample ID	102831344			
Filename	U10721B_10			
Injected By	BAL			
Total Amount Extracted	1057.4 mL	Matrix	WATER	
% Moisture	NA	Dilution	NA	
Dry Weight Extracted	NA	Collected	06/27/2001	
ICAL Date	07/21/2001	Received	06/29/2001	
CCal Filename(s)	U10721B_01 & U10721B_16	Extracted	07/10/2001	
Method Blank ID	BLANK-1153	Analyzed	07/22/2001 03:50	

Native Isomers	Conc ng/L	EMPC ng/L	LRL ng/L		Internal Standards	ng's Added	Percent Recovery
2,3,7,8-TCDF	ND	----	0.0052	A	2,3,7,8-TCDF-13C	2.00	84
Total TCDF	ND	----	0.0019		2,3,7,8-TCDD-13C	2.00	77
					1,2,3,7,8-PeCDF-13C	2.00	94
2,3,7,8-TCDD	ND	----	0.0081	A	2,3,4,7,8-PeCDF-13C	2.00	106
Total TCDD	ND	----	0.0019		1,2,3,7,8-PeCDD-13C	2.00	96
					1,2,3,4,7,8-HxCDF-13C	2.00	110
1,2,3,7,8-PeCDF	ND	----	0.0095		1,2,3,6,7,8-HxCDF-13C	2.00	129
2,3,4,7,8-PeCDF	ND	----	0.0095		2,3,4,6,7,8-HxCDF-13C	2.00	125
Total PeCDF	ND	----	0.0095		1,2,3,7,8,9-HxCDF-13C	2.00	104
					1,2,3,4,7,8-HxCDD-13C	2.00	149
1,2,3,7,8-PeCDD	ND	----	0.0095		1,2,3,6,7,8-HxCDD-13C	2.00	131
Total PeCDD	ND	----	0.0095		1,2,3,4,6,7,8-HpCDF-13C	2.00	97
					1,2,3,4,7,8,9-HpCDF-13C	2.00	85
1,2,3,4,7,8-HxCDF	ND	----	0.0095		1,2,3,4,6,7,8-HpCDD-13C	2.00	86
1,2,3,6,7,8-HxCDF	ND	----	0.0095		OCDD-13C	4.00	77
2,3,4,6,7,8-HxCDF	ND	----	0.0095				
1,2,3,7,8,9-HxCDF	ND	----	0.0095		1,2,3,4-TCDD-13C	2.00	NA
Total HxCDF	ND	----	0.0095		1,2,3,7,8,9-HxCDD-13C	2.00	NA
1,2,3,4,7,8-HxCDD	ND	----	0.0095		2,3,7,8-TCDD-37Cl4	0.20	86
1,2,3,6,7,8-HxCDD	ND	----	0.0095				
1,2,3,7,8,9-HxCDD	ND	----	0.0095				
Total HxCDD	ND	----	0.0095				
1,2,3,4,6,7,8-HpCDF	----	0.45	0.0120	EA	Total 2,3,7,8-TCDD		
1,2,3,4,7,8,9-HpCDF	ND	----	0.0160	A	Equivalence: 0.0043 ng/L		
Total HpCDF	ND	----	0.0095		(Using ITE Factors)		
1,2,3,4,6,7,8-HpCDD	0.20	----	0.0160	A			
Total HpCDD	0.20	----	0.0095				
OCDF	0.24	----	0.0390	A			
OCDD	2.10	----	0.0560	A			

Conc = Concentration (Totals include 2,3,7,8-substituted isomers)
 EMPC = Estimated Maximum Possible Concentration
 A = Detection Limit based on signal-to-noise measurement
 J = Concentration detected is below the calibration range
 B = Less than 10 times higher than method blank level
 P = Recovery outside of target range
 Nn = Value obtained from additional analysis

LRL = Lower Reporting Limit
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Report No.....01-1046199

REPORT OF LABORATORY ANALYSIS

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Method 8290 Analysis Results

Client - ROBERT E LEE

Client's Sample ID	01-10546		
Lab Sample ID	102831351		
Filename	U10721B_11		
Injected By	BAL		
Total Amount Extracted	1051.8 mL	Matrix	WATER
% Moisture	NA	Dilution	NA
Dry Weight Extracted	NA	Collected	06/27/2001
ICAL Date	07/21/2001	Received	06/29/2001
CCal Filename(s)	U10721B_01 & U10721B_16	Extracted	07/10/2001
Method Blank ID	BLANK-1153	Analyzed	07/22/2001 04:40

Native Isomers	Conc ng/L	EMPC ng/L	LRL ng/L		Internal Standards	ng's Added	Percent Recovery
2,3,7,8-TCDF	ND	-----	0.0029	A	2,3,7,8-TCDF-13C	2.00	89
Total TCDF	0.0092	-----	0.0019	J	2,3,7,8-TCDD-13C	2.00	82
					1,2,3,7,8-PeCDF-13C	2.00	100
2,3,7,8-TCDD	ND	-----	0.0064	A	2,3,4,7,8-PeCDF-13C	2.00	117
Total TCDD	ND	-----	0.0019		1,2,3,7,8-PeCDD-13C	2.00	106
					1,2,3,4,7,8-HxCDF-13C	2.00	115
1,2,3,7,8-PeCDF	ND	-----	0.0095		1,2,3,6,7,8-HxCDF-13C	2.00	114
2,3,4,7,8-PeCDF	ND	-----	0.0095		2,3,4,6,7,8-HxCDF-13C	2.00	125
Total PeCDF	ND	-----	0.0095		1,2,3,7,8,9-HxCDF-13C	2.00	101
					1,2,3,4,7,8-HxCDD-13C	2.00	122
1,2,3,7,8-PeCDD	ND	-----	0.0095		1,2,3,6,7,8-HxCDD-13C	2.00	119
Total PeCDD	ND	-----	0.0095		1,2,3,4,6,7,8-HpCDF-13C	2.00	87
					1,2,3,4,7,8,9-HpCDF-13C	2.00	71
1,2,3,4,7,8-HxCDF	ND	-----	0.0095		1,2,3,4,6,7,8-HpCDD-13C	2.00	76
1,2,3,6,7,8-HxCDF	ND	-----	0.0100	A	OCDD-13C	4.00	66
2,3,4,6,7,8-HxCDF	ND	-----	0.0095				
1,2,3,7,8,9-HxCDF	ND	-----	0.0095		1,2,3,4-TCDD-13C	2.00	NA
Total HxCDF	0.0750	-----	0.0095		1,2,3,7,8,9-HxCDD-13C	2.00	NA
1,2,3,4,7,8-HxCDD	ND	-----	0.0095		2,3,7,8-TCDD-37Cl4	0.20	88
1,2,3,6,7,8-HxCDD	-----	0.021	0.0095	I			
1,2,3,7,8,9-HxCDD	ND	-----	0.0095				
Total HxCDD	0.0460	-----	0.0095	J			
1,2,3,4,6,7,8-HpCDF	-----	0.720	0.0095	E	Total 2,3,7,8-TCDD		
1,2,3,4,7,8,9-HpCDF	ND	-----	0.0140	A	Equivalence: 0.013 ng/L		
Total HpCDF	ND	-----	0.0095		(Using ITE Factors)		
1,2,3,4,6,7,8-HpCDD	0.6100	-----	0.0210	A			
Total HpCDD	1.2000	-----	0.0095				
OCDF	0.5100	-----	0.0310	A			
OCDD	6.6000	-----	0.0760	A			

Conc = Concentration (Totals include 2,3,7,8-substituted isomers)
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Report No.....01-1046199

REPORT OF LABORATORY ANALYSIS

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Method 8290 Analysis Results

Client - ROBERT E LEE

Client's Sample ID	01-10547		
Lab Sample ID	102831369		
Filename	U10721B_12		
Injected By	BAL		
Total Amount Extracted	1049.6 mL	Matrix	WATER
% Moisture	NA	Dilution	NA
Dry Weight Extracted	NA	Collected	06/27/2001
ICAL Date	07/21/2001	Received	06/29/2001
CCal Filename(s)	U10721B_01 & U10721B_16	Extracted	07/10/2001
Method Blank ID	BLANK-1153	Analyzed	07/22/2001 05:30

Native Isomers	Conc ng/L	EMPC ng/L	LRL ng/L		Internal Standards	ng's Added	Percent Recovery
2,3,7,8-TCDF	0.0056	-----	0.0022	JA	2,3,7,8-TCDF-13C	2.00	86
Total TCDF	0.0160	-----	0.0019		2,3,7,8-TCDD-13C	2.00	80
					1,2,3,7,8-PeCDF-13C	2.00	85
2,3,7,8-TCDD	ND	-----	0.0053	A	2,3,4,7,8-PeCDF-13C	2.00	88
Total TCDD	ND	-----	0.0019		1,2,3,7,8-PeCDD-13C	2.00	79
					1,2,3,4,7,8-HxCDF-13C	2.00	106
1,2,3,7,8-PeCDF	-----	0.015	0.0110	IA	1,2,3,6,7,8-HxCDF-13C	2.00	122
2,3,4,7,8-PeCDF	0.0210	-----	0.0095	J	2,3,4,6,7,8-HxCDF-13C	2.00	115
Total PeCDF	0.0870	-----	0.0095		1,2,3,7,8,9-HxCDF-13C	2.00	101
					1,2,3,4,7,8-HxCDD-13C	2.00	132
1,2,3,7,8-PeCDD	ND	-----	0.0095		1,2,3,6,7,8-HxCDD-13C	2.00	118
Total PeCDD	ND	-----	0.0095		1,2,3,4,6,7,8-HpCDF-13C	2.00	89
					1,2,3,4,7,8,9-HpCDF-13C	2.00	71
1,2,3,4,7,8-HxCDF	0.0350	-----	0.0095	J	1,2,3,4,6,7,8-HpCDD-13C	2.00	75
1,2,3,6,7,8-HxCDF	ND	-----	0.0095		OCDD-13C	4.00	66
2,3,4,6,7,8-HxCDF	0.0210	-----	0.0095	J			
1,2,3,7,8,9-HxCDF	0.0220	-----	0.0095	J	1,2,3,4-TCDD-13C	2.00	NA
Total HxCDF	0.7100	-----	0.0095		1,2,3,7,8,9-HxCDD-13C	2.00	NA
1,2,3,4,7,8-HxCDD	ND	-----	0.0095		2,3,7,8-TCDD-37Cl4	0.20	87
1,2,3,6,7,8-HxCDD	0.2600	-----	0.0095				
1,2,3,7,8,9-HxCDD	0.0250	-----	0.0095	J			
Total HxCDD	0.7700	-----	0.0095				
1,2,3,4,6,7,8-HpCDF	-----	0.730	0.0095	E	Total 2,3,7,8-TCDD		
1,2,3,4,7,8,9-HpCDF	ND	-----	0.0140	A	Equivalence: 0.12 ng/L		
Total HpCDF	0.6500	-----	0.0095		(Using ITE Factors)		
1,2,3,4,6,7,8-HpCDD	4.2000	-----	0.0280	A			
Total HpCDD	7.8000	-----	0.0095				
OCDF	0.3800	-----	0.0530	A			
OCDD	26.0000	-----	0.0500	A			

Conc = Concentration (Totals include 2,3,7,8-substituted isomers)
EMPC = Estimated Maximum Possible Concentration
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Report No.....01-1046199

REPORT OF LABORATORY ANALYSIS

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Method 8290 Analysis Results

Client - ROBERT E LEE

Client's Sample ID	01-10548		
Lab Sample ID	102831377		
Filename	U10721B_13		
Injected By	BAL		
Total Amount Extracted	1055.4 mL	Matrix	WATER
% Moisture	NA	Dilution	NA
Dry Weight Extracted	NA	Collected	06/27/2001
ICAL Date	07/21/2001	Received	06/29/2001
CCal Filename(s)	U10721B_01 & U10721B_16	Extracted	07/10/2001
Method Blank ID	BLANK-1153	Analyzed	07/22/2001 06:20

Native Isomers	Conc ng/L	EMPC ng/L	LRL ng/L		Internal Standards	ng's Added	Percent Recovery
2,3,7,8-TCDF	0.170	-----	0.00190		2,3,7,8-TCDF-13C	2.00	89
Total TCDF	1.500	-----	0.00190		2,3,7,8-TCDD-13C	2.00	96
					1,2,3,7,8-PeCDF-13C	2.00	103
2,3,7,8-TCDD	ND	-----	0.00280	A	2,3,4,7,8-PeCDF-13C	2.00	71
Total TCDD	ND	-----	0.00190		1,2,3,7,8-PeCDD-13C	2.00	66
					1,2,3,4,7,8-HxCDF-13C	2.00	111
1,2,3,7,8-PeCDF	0.660	-----	0.00950		1,2,3,6,7,8-HxCDF-13C	2.00	155 IP
2,3,4,7,8-PeCDF	0.690	-----	0.00950		2,3,4,6,7,8-HxCDF-13C	2.00	110
Total PeCDF	2.700	-----	0.00950		1,2,3,7,8,9-HxCDF-13C	2.00	106
					1,2,3,4,7,8-HxCDD-13C	2.00	101
1,2,3,7,8-PeCDD	0.049	-----	0.00950		1,2,3,6,7,8-HxCDD-13C	2.00	100
Total PeCDD	0.150	-----	0.00950		1,2,3,4,6,7,8-HpCDF-13C	2.00	113 I
					1,2,3,4,7,8,9-HpCDF-13C	2.00	59
1,2,3,4,7,8-HxCDF	2.200	-----	0.01900	A	1,2,3,4,6,7,8-HpCDD-13C	2.00	100 I
1,2,3,6,7,8-HxCDF	7.800	-----	0.01400	A	OCDD-13C	4.00	221 IP
2,3,4,6,7,8-HxCDF	2.300	-----	0.01300	A			
1,2,3,7,8,9-HxCDF	1.600	-----	0.00990	A	1,2,3,4-TCDD-13C	2.00	NA
Total HxCDF	97.000	-----	0.00950		1,2,3,7,8,9-HxCDD-13C	2.00	NA
1,2,3,4,7,8-HxCDD	0.220	-----	0.01300	A	2,3,7,8-TCDD-37Cl4	0.20	101
1,2,3,6,7,8-HxCDD	21.000	-----	0.01900	A			
1,2,3,7,8,9-HxCDD	1.100	-----	0.01700	A			
Total HxCDD	55.000	-----	0.00950				
1,2,3,4,6,7,8-HpCDF	-----	290	0.03000	EA	Total 2,3,7,8-TCDD		
1,2,3,4,7,8,9-HpCDF	4.000	-----	0.03200	A	Equivalence: 8.5 ng/L		
Total HpCDF	180.000	-----	0.00950		(Using ITE Factors)		
1,2,3,4,6,7,8-HpCDD	300.000	-----	0.01300	A			
Total HpCDD	520.000	-----	0.00950				
OCDF	110.000	-----	0.01900				
OCDD	1300.000	-----	0.01900				

Conc = Concentration (Totals include 2,3,7,8-substituted isomers)
EMPC = Estimated Maximum Possible Concentration
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Report No.....01-1046199

REPORT OF LABORATORY ANALYSIS

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Method 8290 Analysis Results

Client - ROBERT E LEE

Client's Sample ID	01-10549		
Lab Sample ID	102831385		
Filename	U10721B_14		
Injected By	BAL		
Total Amount Extracted	1054.9 mL	Matrix	WATER
% Moisture	NA	Dilution	NA
Dry Weight Extracted	NA	Collected	06/27/2001
ICAL Date	07/21/2001	Received	06/29/2001
CCal Filename(s)	U10721B_01 & U10721B_16	Extracted	07/10/2001
Method Blank ID	BLANK-1153	Analyzed	07/22/2001 07:10

Native Isomers	Conc ng/L	EMPC ng/L	LRL ng/L		Internal Standards	ng's Added	Percent Recovery
2,3,7,8-TCDF	ND	----	0.0053	A	2,3,7,8-TCDF-13C	2.00	77
Total TCDF	ND	----	0.0019		2,3,7,8-TCDD-13C	2.00	71
					1,2,3,7,8-PeCDF-13C	2.00	90
2,3,7,8-TCDD	ND	----	0.0089	A	2,3,4,7,8-PeCDF-13C	2.00	106
Total TCDD	ND	----	0.0019		1,2,3,7,8-PeCDD-13C	2.00	91
					1,2,3,4,7,8-HxCDF-13C	2.00	110
1,2,3,7,8-PeCDF	ND	----	0.0095		1,2,3,6,7,8-HxCDF-13C	2.00	114
2,3,4,7,8-PeCDF	ND	----	0.0095		2,3,4,6,7,8-HxCDF-13C	2.00	124
Total PeCDF	ND	----	0.0095		1,2,3,7,8,9-HxCDF-13C	2.00	99
					1,2,3,4,7,8-HxCDD-13C	2.00	130
1,2,3,7,8-PeCDD	ND	----	0.0095		1,2,3,6,7,8-HxCDD-13C	2.00	128
Total PeCDD	ND	----	0.0095		1,2,3,4,6,7,8-HpCDF-13C	2.00	90
					1,2,3,4,7,8,9-HpCDF-13C	2.00	74
1,2,3,4,7,8-HxCDF	ND	----	0.0095		1,2,3,4,6,7,8-HpCDD-13C	2.00	78
1,2,3,6,7,8-HxCDF	ND	----	0.0095		OCDD-13C	4.00	70
2,3,4,6,7,8-HxCDF	ND	----	0.0095				
1,2,3,7,8,9-HxCDF	ND	----	0.0110	A	1,2,3,4-TCDD-13C	2.00	NA
Total HxCDF	0.089	----	0.0095		1,2,3,7,8,9-HxCDD-13C	2.00	NA
1,2,3,4,7,8-HxCDD	ND	----	0.0095		2,3,7,8-TCDD-37Cl4	0.20	75
1,2,3,6,7,8-HxCDD	ND	----	0.0095				
1,2,3,7,8,9-HxCDD	ND	----	0.0095				
Total HxCDD	ND	----	0.0095				
1,2,3,4,6,7,8-HpCDF	----	0.17	0.0150	EA	Total 2,3,7,8-TCDD		
1,2,3,4,7,8,9-HpCDF	ND	----	0.0270	A	Equivalence: 0.0040 ng/L		
Total HpCDF	ND	----	0.0095		(Using ITE Factors)		
1,2,3,4,6,7,8-HpCDD	0.150	----	0.0250	A			
Total HpCDD	0.240	----	0.0095				
OCDF	0.280	----	0.0450	A			
OCDD	2.300	----	0.0760	A			

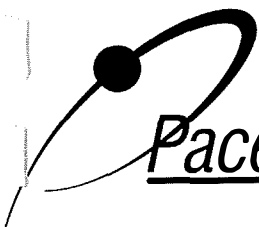
Conc = Concentration (Totals include 2,3,7,8-substituted isomers)
EMPC = Estimated Maximum Possible Concentration
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REPORT OF LABORATORY ANALYSIS

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Method 8290 Analysis Results

Client - ROBERT E LEE

Client's Sample ID	01-10550		
Lab Sample ID	102831393		
Filename	U10724A_13		
Injected By	CVS		
Total Amount Extracted	1050.3 mL	Matrix	WATER
% Moisture	NA	Dilution	NA
Dry Weight Extracted	NA	Collected	06/27/2001
ICAL Date	07/24/2001	Received	06/29/2001
CCal Filename(s)	U10724A_06 & U10724A_21	Extracted	07/10/2001
Method Blank ID	BLANK-1153	Analyzed	07/24/2001 22:51

Native Isomers	Conc ng/L	EMPC ng/L	LRL ng/L		Internal Standards	ng's Added	Percent Recovery
2,3,7,8-TCDF	ND	-----	0.0088	AN2	2,3,7,8-TCDF-13C	2.00	69 N2
Total TCDF	ND	-----	0.0019	N2	2,3,7,8-TCDD-13C	2.00	67 N2
					1,2,3,7,8-PeCDF-13C	2.00	97
2,3,7,8-TCDD	ND	-----	0.0079	AN2	2,3,4,7,8-PeCDF-13C	2.00	90
Total TCDD	ND	-----	0.0019	N2	1,2,3,7,8-PeCDD-13C	2.00	94
					1,2,3,4,7,8-HxCDF-13C	2.00	94
1,2,3,7,8-PeCDF	ND	-----	0.0220	A	1,2,3,6,7,8-HxCDF-13C	2.00	87
2,3,4,7,8-PeCDF	ND	-----	0.0100	A	2,3,4,6,7,8-HxCDF-13C	2.00	100
Total PeCDF	ND	-----	0.0095		1,2,3,7,8,9-HxCDF-13C	2.00	91
					1,2,3,4,7,8-HxCDD-13C	2.00	95
1,2,3,7,8-PeCDD	ND	-----	0.0310	A	1,2,3,6,7,8-HxCDD-13C	2.00	94
Total PeCDD	ND	-----	0.0095		1,2,3,4,6,7,8-HpCDF-13C	2.00	95
					1,2,3,4,7,8,9-HpCDF-13C	2.00	94
1,2,3,4,7,8-HxCDF	ND	-----	0.0210	A	1,2,3,4,6,7,8-HpCDD-13C	2.00	104
1,2,3,6,7,8-HxCDF	ND	-----	0.0240	A	OCDD-13C	4.00	97
2,3,4,6,7,8-HxCDF	ND	-----	0.0240	A			
1,2,3,7,8,9-HxCDF	ND	-----	0.0370	A	1,2,3,4-TCDD-13C	2.00	NA
Total HxCDF	0.190	-----	0.0095		1,2,3,7,8,9-HxCDD-13C	2.00	NA
1,2,3,4,7,8-HxCDD	ND	-----	0.0360	A	2,3,7,8-TCDD-37Cl4	0.20	61
1,2,3,6,7,8-HxCDD	0.064	-----	0.0450	A			
1,2,3,7,8,9-HxCDD	ND	-----	0.0370	A			
Total HxCDD	0.130	-----	0.0095				
1,2,3,4,6,7,8-HpCDF	-----	0.76	0.0330	EA	Total 2,3,7,8-TCDD		
1,2,3,4,7,8,9-HpCDF	ND	-----	0.0170	A	Equivalence: 0.022 ng/L		
Total HpCDF	0.460	-----	0.0095		(Using ITE Factors)		
1,2,3,4,6,7,8-HpCDD	0.810	-----	0.0360	A			
Total HpCDD	1.300	-----	0.0095				
OCDF	0.550	-----	0.0400	A			
OCDD	6.700	-----	0.0400	A			

Conc = Concentration (Totals include 2,3,7,8-substituted isomers)
EMPC = Estimated Maximum Possible Concentration
A = Detection Limit based on signal-to-noise measurement
J = Concentration detected is below the calibration range
B = Less than 10 times higher than method blank level
P = Recovery outside of target range
Nn = Value obtained from additional analysis

LRL = Lower Reporting Limit
I = Interference
E = PCDE Interference
S = Saturated signal
ND = Not Detected
NA = Not Applicable
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Report No.....01-1046199

REPORT OF LABORATORY ANALYSIS

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Method 8290 Analysis Results

Client - ROBERT E LEE

Client's Sample ID	01-10551		
Lab Sample ID	102831401		
Filename	U10724A_14		
Injected By	CVS		
Total Amount Extracted	1058.7 mL	Matrix	WATER
% Moisture	NA	Dilution	NA
Dry Weight Extracted	NA	Collected	06/27/2001
ICAL Date	07/24/2001	Received	06/29/2001
CCal Filename(s)	U10724A_06 & U10724A_21	Extracted	07/10/2001
Method Blank ID	BLANK-1153	Analyzed	07/24/2001 23:41

Native Isomers	Conc ng/L	EMPC ng/L	LRL ng/L	Internal Standards	ng's Added	Percent Recovery
2,3,7,8-TCDF	ND	-----	0.0047 AN2	2,3,7,8-TCDF-13C	2.00	75 N2
Total TCDF	ND	-----	0.0019 N2	2,3,7,8-TCDD-13C	2.00	72 N2
				1,2,3,7,8-PeCDF-13C	2.00	100
2,3,7,8-TCDD	ND	-----	0.0047 AN2	2,3,4,7,8-PeCDF-13C	2.00	96
Total TCDD	ND	-----	0.0019 N2	1,2,3,7,8-PeCDD-13C	2.00	96
				1,2,3,4,7,8-HxCDF-13C	2.00	99
1,2,3,7,8-PeCDF	ND	-----	0.0190 A	1,2,3,6,7,8-HxCDF-13C	2.00	91
2,3,4,7,8-PeCDF	ND	-----	0.0120 A	2,3,4,6,7,8-HxCDF-13C	2.00	106
Total PeCDF	ND	-----	0.0094	1,2,3,7,8,9-HxCDF-13C	2.00	98
				1,2,3,4,7,8-HxCDD-13C	2.00	101
1,2,3,7,8-PeCDD	ND	-----	0.0320 A	1,2,3,6,7,8-HxCDD-13C	2.00	97
Total PeCDD	ND	-----	0.0094	1,2,3,4,6,7,8-HpCDF-13C	2.00	95
				1,2,3,4,7,8,9-HpCDF-13C	2.00	94
1,2,3,4,7,8-HxCDF	ND	-----	0.0096 A	1,2,3,4,6,7,8-HpCDD-13C	2.00	99
1,2,3,6,7,8-HxCDF	ND	-----	0.0094	OCDD-13C	4.00	89
2,3,4,6,7,8-HxCDF	ND	-----	0.0130 A			
1,2,3,7,8,9-HxCDF	ND	-----	0.0290 A	1,2,3,4-TCDD-13C	2.00	NA
Total HxCDF	0.023	-----	0.0094 J	1,2,3,7,8,9-HxCDD-13C	2.00	NA
1,2,3,4,7,8-HxCDD	ND	-----	0.0360 A	2,3,7,8-TCDD-37Cl4	0.20	74
1,2,3,6,7,8-HxCDD	ND	-----	0.0300 A			
1,2,3,7,8,9-HxCDD	ND	-----	0.0410 A			
Total HxCDD	ND	-----	0.0094			
1,2,3,4,6,7,8-HpCDF	-----	0.19	0.0210 EIA	Total 2,3,7,8-TCDD		
1,2,3,4,7,8,9-HpCDF	ND	-----	0.0260 A	Equivalence: 0.0046 ng/L		
Total HpCDF	0.088	-----	0.0094	(Using ITE Factors)		
1,2,3,4,6,7,8-HpCDD	0.290	-----	0.0460 A			
Total HpCDD	0.290	-----	0.0094			
OCDF	0.130	-----	0.0190			
OCDD	1.600	-----	0.0510 A			

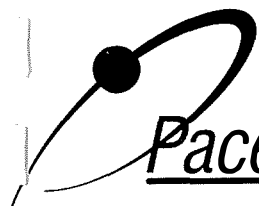
Conc = Concentration (Totals include 2,3,7,8-substituted isomers)
EMPC = Estimated Maximum Possible Concentration
A = Detection Limit based on signal-to-noise measurement
J = Concentration detected is below the calibration range
B = Less than 10 times higher than method blank level
P = Recovery outside of target range
Nn = Value obtained from additional analysis

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I = Interference
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Report No.....01-1046199

REPORT OF LABORATORY ANALYSIS

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Method 8290 Analysis Results

Client - ROBERT E LEE

Client's Sample ID	01-10552		
Lab Sample ID	102831419		
Filename	U10724A_15		
Injected By	CVS		
Total Amount Extracted	1055.6 mL	Matrix	WATER
% Moisture	NA	Dilution	NA
Dry Weight Extracted	NA	Collected	06/27/2001
ICAL Date	07/24/2001	Received	06/29/2001
CCal Filename(s)	U10724A_06 & U10724A_21	Extracted	07/10/2001
Method Blank ID	BLANK-1153	Analyzed	07/25/2001 00:31

Native Isomers	Conc ng/L	EMPC ng/L	LRL ng/L		Internal Standards	ng's Added	Percent Recovery
2,3,7,8-TCDF	ND	-----	0.0048	AN2	2,3,7,8-TCDF-13C	2.00	82 N2
Total TCDF	ND	-----	0.0019	N2	2,3,7,8-TCDD-13C	2.00	76 N2
					1,2,3,7,8-PeCDF-13C	2.00	106
2,3,7,8-TCDD	ND	-----	0.0074	AN2	2,3,4,7,8-PeCDF-13C	2.00	99
Total TCDD	ND	-----	0.0019	N2	1,2,3,7,8-PeCDD-13C	2.00	102
					1,2,3,4,7,8-HxCDF-13C	2.00	113
1,2,3,7,8-PeCDF	ND	-----	0.0190	A	1,2,3,6,7,8-HxCDF-13C	2.00	98
2,3,4,7,8-PeCDF	ND	-----	0.0200	A	2,3,4,6,7,8-HxCDF-13C	2.00	112
Total PeCDF	ND	-----	0.0095		1,2,3,7,8,9-HxCDF-13C	2.00	100
					1,2,3,4,7,8-HxCDD-13C	2.00	105
1,2,3,7,8-PeCDD	ND	-----	0.0250	A	1,2,3,6,7,8-HxCDD-13C	2.00	99
Total PeCDD	ND	-----	0.0095		1,2,3,4,6,7,8-HpCDF-13C	2.00	105
					1,2,3,4,7,8,9-HpCDF-13C	2.00	98
1,2,3,4,7,8-HxCDF	ND	-----	0.0210	A	1,2,3,4,6,7,8-HpCDD-13C	2.00	111
1,2,3,6,7,8-HxCDF	ND	-----	0.0200	A	OCDD-13C	4.00	91
2,3,4,6,7,8-HxCDF	ND	-----	0.0150	A			
1,2,3,7,8,9-HxCDF	ND	-----	0.0310	A	1,2,3,4-TCDD-13C	2.00	NA
Total HxCDF	ND	-----	0.0095		1,2,3,7,8,9-HxCDD-13C	2.00	NA
1,2,3,4,7,8-HxCDD	ND	-----	0.0330	A	2,3,7,8-TCDD-37Cl4	0.20	73
1,2,3,6,7,8-HxCDD	ND	-----	0.0430	A			
1,2,3,7,8,9-HxCDD	ND	-----	0.0360	A			
Total HxCDD	ND	-----	0.0095				
1,2,3,4,6,7,8-HpCDF	-----	0.035	0.0220	IA	Total 2,3,7,8-TCDD		
1,2,3,4,7,8,9-HpCDF	ND	-----	0.0350	A	Equivalence: 0.0015 ng/L		
Total HpCDF	ND	-----	0.0095		(Using ITE Factors)		
1,2,3,4,6,7,8-HpCDD	0.090	-----	0.0330	A			
Total HpCDD	0.090	-----	0.0095				
OCDF	-----	0.038	0.0370	IA			
OCDD	0.580	-----	0.0480	A			

Conc = Concentration (Totals include 2,3,7,8-substituted isomers)
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B = Less than 10 times higher than method blank level
P = Recovery outside of target range
Nn = Value obtained from additional analysis

LRL = Lower Reporting Limit
I = Interference
E = PCDE Interference
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Report No.....01-1046199

REPORT OF LABORATORY ANALYSIS

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Method 8290 Analysis Results

Client - ROBERT E LEE

Client's Sample ID	01-10553		
Lab Sample ID	102831427		
Filename	U10724A_16		
Injected By	CVS		
Total Amount Extracted	1051.1 mL	Matrix	WATER
% Moisture	NA	Dilution	NA
Dry Weight Extracted	NA	Collected	06/27/2001
ICAL Date	07/24/2001	Received	06/29/2001
CCal Filename(s)	U10724A_06 & U10724A_21	Extracted	07/10/2001
Method Blank ID	BLANK-1153	Analyzed	07/25/2001 01:21

Native Isomers	Conc ng/L	EMPC ng/L	LRL ng/L	Internal Standards	ng's Added	Percent Recovery
2,3,7,8-TCDF	ND	-----	0.0049 AN2	2,3,7,8-TCDF-13C	2.00	73 N2
Total TCDF	ND	-----	0.0019 N2	2,3,7,8-TCDD-13C	2.00	68 N2
				1,2,3,7,8-PeCDF-13C	2.00	92
2,3,7,8-TCDD	ND	-----	0.0056 AN2	2,3,4,7,8-PeCDF-13C	2.00	88
Total TCDD	ND	-----	0.0019 N2	1,2,3,7,8-PeCDD-13C	2.00	89
				1,2,3,4,7,8-HxCDF-13C	2.00	95
1,2,3,7,8-PeCDF	ND	-----	0.0230 A	1,2,3,6,7,8-HxCDF-13C	2.00	86
2,3,4,7,8-PeCDF	ND	-----	0.0170 A	2,3,4,6,7,8-HxCDF-13C	2.00	96
Total PeCDF	ND	-----	0.0095	1,2,3,7,8,9-HxCDF-13C	2.00	94
				1,2,3,4,7,8-HxCDD-13C	2.00	94
1,2,3,7,8-PeCDD	ND	-----	0.0450 A	1,2,3,6,7,8-HxCDD-13C	2.00	87
Total PeCDD	ND	-----	0.0095	1,2,3,4,6,7,8-HpCDF-13C	2.00	86
				1,2,3,4,7,8,9-HpCDF-13C	2.00	91
1,2,3,4,7,8-HxCDF	ND	-----	0.0270 A	1,2,3,4,6,7,8-HpCDD-13C	2.00	92
1,2,3,6,7,8-HxCDF	ND	-----	0.0220 A	OCDD-13C	4.00	87
2,3,4,6,7,8-HxCDF	ND	-----	0.0180 A			
1,2,3,7,8,9-HxCDF	ND	-----	0.0160 A	1,2,3,4-TCDD-13C	2.00	NA
Total HxCDF	ND	-----	0.0095	1,2,3,7,8,9-HxCDD-13C	2.00	NA
1,2,3,4,7,8-HxCDD	ND	-----	0.0480 A	2,3,7,8-TCDD-37Cl4	0.20	66
1,2,3,6,7,8-HxCDD	ND	-----	0.0420 A			
1,2,3,7,8,9-HxCDD	ND	-----	0.0350 A			
Total HxCDD	ND	-----	0.0095			
1,2,3,4,6,7,8-HpCDF	ND	-----	0.0240 A	Total 2,3,7,8-TCDD		
1,2,3,4,7,8,9-HpCDF	ND	-----	0.0340 A	Equivalence: 0.00 ng/L		
Total HpCDF	ND	-----	0.0095	(Using ITE Factors)		
1,2,3,4,6,7,8-HpCDD	ND	-----	0.0460 A			
Total HpCDD	ND	-----	0.0095			
OCDF	ND	-----	0.0310 A			
OCDD	ND	-----	0.0630 A			

Conc = Concentration (Totals include 2,3,7,8-substituted isomers)
EMPC = Estimated Maximum Possible Concentration
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B = Less than 10 times higher than method blank level
P = Recovery outside of target range
Nn = Value obtained from additional analysis

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I = Interference
E = PCDE Interference
S = Saturated signal
ND = Not Detected
NA = Not Applicable
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Report No.....01-1046199

REPORT OF LABORATORY ANALYSIS

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Method 8290 Analysis Results

Client - ROBERT E LEE

Client's Sample ID	01-10554		
Lab Sample ID	102831435		
Filename	U10724A_17		
Injected By	CVS		
Total Amount Extracted	1052 mL	Matrix	WATER
% Moisture	NA	Dilution	NA
Dry Weight Extracted	NA	Collected	06/27/2001
ICAL Date	07/24/2001	Received	06/29/2001
CCal Filename(s)	U10724A_06 & U10724A_21	Extracted	07/10/2001
Method Blank ID	BLANK-1153	Analyzed	07/25/2001 02:11

Native Isomers	Conc ng/L	EMPC ng/L	LRL ng/L	Internal Standards	ng's Added	Percent Recovery
2,3,7,8-TCDF	ND	-----	0.0049 AN2	2,3,7,8-TCDF-13C	2.00	73 N2
Total TCDF	ND	-----	0.0019 N2	2,3,7,8-TCDD-13C	2.00	69 N2
				1,2,3,7,8-PeCDF-13C	2.00	96
2,3,7,8-TCDD	ND	-----	0.0084 AN2	2,3,4,7,8-PeCDF-13C	2.00	93
Total TCDD	ND	-----	0.0019 N2	1,2,3,7,8-PeCDD-13C	2.00	95
				1,2,3,4,7,8-HxCDF-13C	2.00	100
1,2,3,7,8-PeCDF	ND	-----	0.0200 A	1,2,3,6,7,8-HxCDF-13C	2.00	88
2,3,4,7,8-PeCDF	ND	-----	0.0180 A	2,3,4,6,7,8-HxCDF-13C	2.00	103
Total PeCDF	ND	-----	0.0095	1,2,3,7,8,9-HxCDF-13C	2.00	95
				1,2,3,4,7,8-HxCDD-13C	2.00	102
1,2,3,7,8-PeCDD	ND	-----	0.0280 A	1,2,3,6,7,8-HxCDD-13C	2.00	94
Total PeCDD	ND	-----	0.0095	1,2,3,4,6,7,8-HpCDF-13C	2.00	97
				1,2,3,4,7,8,9-HpCDF-13C	2.00	92
1,2,3,4,7,8-HxCDF	ND	-----	0.0220 A	1,2,3,4,6,7,8-HpCDD-13C	2.00	96
1,2,3,6,7,8-HxCDF	ND	-----	0.0190 A	OCDD-13C	4.00	95
2,3,4,6,7,8-HxCDF	ND	-----	0.0160 A			
1,2,3,7,8,9-HxCDF	ND	-----	0.0210 A	1,2,3,4-TCDD-13C	2.00	NA
Total HxCDF	ND	-----	0.0095	1,2,3,7,8,9-HxCDD-13C	2.00	NA
1,2,3,4,7,8-HxCDD	ND	-----	0.0320 A	2,3,7,8-TCDD-37Cl4	0.20	69
1,2,3,6,7,8-HxCDD	ND	-----	0.0240 A			
1,2,3,7,8,9-HxCDD	ND	-----	0.0330 A			
Total HxCDD	ND	-----	0.0095			
1,2,3,4,6,7,8-HpCDF	ND	-----	0.0300 A	Total 2,3,7,8-TCDD		
1,2,3,4,7,8,9-HpCDF	ND	-----	0.0340 A	Equivalence: 0.00 ng/L		
Total HpCDF	ND	-----	0.0095	(Using ITE Factors)		
1,2,3,4,6,7,8-HpCDD	ND	-----	0.0330 A			
Total HpCDD	ND	-----	0.0095			
OCDF	ND	-----	0.0330 A			
OCDD	ND	-----	0.0420 A			

Conc = Concentration (Totals include 2,3,7,8-substituted isomers)
EMPC = Estimated Maximum Possible Concentration
A = Detection Limit based on signal-to-noise measurement
J = Concentration detected is below the calibration range
B = Less than 10 times higher than method blank level
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Nn = Value obtained from additional analysis

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Report No.....01-1046199

REPORT OF LABORATORY ANALYSIS

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Method 8290 Analysis Results

Client - ROBERT E LEE

Client's Sample ID	01-10557		
Lab Sample ID	102831443		
Filename	U10724A_18		
Injected By	CVS		
Total Amount Extracted	1058.6 mL	Matrix	WATER
% Moisture	NA	Dilution	NA
Dry Weight Extracted	NA	Collected	06/27/2001
ICAL Date	07/24/2001	Received	06/29/2001
CCal Filename(s)	U10724A_06 & U10724A_21	Extracted	07/10/2001
Method Blank ID	BLANK-1153	Analyzed	07/25/2001 03:01

Native Isomers	Conc ng/L	EMPC ng/L	LRL ng/L		Internal Standards	ng's Added	Percent Recovery
2,3,7,8-TCDF	ND	-----	0.0068	AN2	2,3,7,8-TCDF-13C	2.00	87 N2
Total TCDF	ND	-----	0.0019	N2	2,3,7,8-TCDD-13C	2.00	79 N2
					1,2,3,7,8-PeCDF-13C	2.00	102
2,3,7,8-TCDD	ND	-----	0.0054	AN2	2,3,4,7,8-PeCDF-13C	2.00	96
Total TCDD	ND	-----	0.0019	N2	1,2,3,7,8-PeCDD-13C	2.00	101
					1,2,3,4,7,8-HxCDF-13C	2.00	104
1,2,3,7,8-PeCDF	ND	-----	0.0130	A	1,2,3,6,7,8-HxCDF-13C	2.00	98
2,3,4,7,8-PeCDF	ND	-----	0.0140	A	2,3,4,6,7,8-HxCDF-13C	2.00	111
Total PeCDF	0.026	-----	0.0094	J	1,2,3,7,8,9-HxCDF-13C	2.00	107
					1,2,3,4,7,8-HxCDD-13C	2.00	111
1,2,3,7,8-PeCDD	ND	-----	0.0130	A	1,2,3,6,7,8-HxCDD-13C	2.00	99
Total PeCDD	ND	-----	0.0094		1,2,3,4,6,7,8-HpCDF-13C	2.00	108
					1,2,3,4,7,8,9-HpCDF-13C	2.00	107
1,2,3,4,7,8-HxCDF	-----	0.010	0.0094	I	1,2,3,4,6,7,8-HpCDD-13C	2.00	113
1,2,3,6,7,8-HxCDF	-----	0.010	0.0110	IA	OCDD-13C	4.00	125
2,3,4,6,7,8-HxCDF	ND	-----	0.0110	A			
1,2,3,7,8,9-HxCDF	ND	-----	0.0110	A	1,2,3,4-TCDD-13C	2.00	NA
Total HxCDF	0.280	-----	0.0094		1,2,3,7,8,9-HxCDD-13C	2.00	NA
1,2,3,4,7,8-HxCDD	ND	-----	0.0095	A	2,3,7,8-TCDD-37Cl4	0.20	76
1,2,3,6,7,8-HxCDD	0.093	-----	0.0240	A			
1,2,3,7,8,9-HxCDD	ND	-----	0.0220	A			
Total HxCDD	0.200	-----	0.0094				
1,2,3,4,6,7,8-HpCDF	-----	1.300	0.0210	EA	Total 2,3,7,8-TCDD		
1,2,3,4,7,8,9-HpCDF	ND	-----	0.0290	A	Equivalence: 0.047 ng/L		
Total HpCDF	0.910	-----	0.0094		(Using ITE Factors)		
1,2,3,4,6,7,8-HpCDD	1.800	-----	0.0380	A			
Total HpCDD	2.900	-----	0.0094				
OCDF	1.400	-----	0.0190				
OCDD	19.000	-----	0.0190				

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Method 8290 Analysis Results

Client - ROBERT E LEE

Client's Sample ID	01-10560		
Lab Sample ID	102831450		
Filename	U10724A_19		
Injected By	CVS		
Total Amount Extracted	1059.7 mL	Matrix	WATER
% Moisture	NA	Dilution	NA
Dry Weight Extracted	NA	Collected	06/27/2001
ICAL Date	07/24/2001	Received	06/29/2001
CCal Filename(s)	U10724A_06 & U10724A_21	Extracted	07/10/2001
Method Blank ID	BLANK-1153	Analyzed	07/25/2001 03:51

Native Isomers	Conc ng/L	EMPC ng/L	LRL ng/L		Internal Standards	ng's Added	Percent Recovery
2,3,7,8-TCDF	ND	-----	0.0050	AN2	2,3,7,8-TCDF-13C	2.00	78 N2
Total TCDF	ND	-----	0.0019	N2	2,3,7,8-TCDD-13C	2.00	72 N2
					1,2,3,7,8-PeCDF-13C	2.00	98
2,3,7,8-TCDD	ND	-----	0.0067	AN2	2,3,4,7,8-PeCDF-13C	2.00	93
Total TCDD	ND	-----	0.0019	N2	1,2,3,7,8-PeCDD-13C	2.00	99
					1,2,3,4,7,8-HxCDF-13C	2.00	95
1,2,3,7,8-PeCDF	ND	-----	0.0140	A	1,2,3,6,7,8-HxCDF-13C	2.00	85
2,3,4,7,8-PeCDF	ND	-----	0.0096	A	2,3,4,6,7,8-HxCDF-13C	2.00	103
Total PeCDF	ND	-----	0.0094		1,2,3,7,8,9-HxCDF-13C	2.00	97
					1,2,3,4,7,8-HxCDD-13C	2.00	104
1,2,3,7,8-PeCDD	ND	-----	0.0170	A	1,2,3,6,7,8-HxCDD-13C	2.00	92
Total PeCDD	ND	-----	0.0094		1,2,3,4,6,7,8-HpCDF-13C	2.00	94
					1,2,3,4,7,8,9-HpCDF-13C	2.00	98
1,2,3,4,7,8-HxCDF	ND	-----	0.0100	A	1,2,3,4,6,7,8-HpCDD-13C	2.00	103
1,2,3,6,7,8-HxCDF	ND	-----	0.0110	A	OCDD-13C	4.00	90
2,3,4,6,7,8-HxCDF	ND	-----	0.0097	A			
1,2,3,7,8,9-HxCDF	ND	-----	0.0130	A	1,2,3,4-TCDD-13C	2.00	NA
Total HxCDF	ND	-----	0.0094		1,2,3,7,8,9-HxCDD-13C	2.00	NA
1,2,3,4,7,8-HxCDD	ND	-----	0.0150	A	2,3,7,8-TCDD-37Cl4	0.20	76
1,2,3,6,7,8-HxCDD	ND	-----	0.0170	A			
1,2,3,7,8,9-HxCDD	ND	-----	0.0150	A			
Total HxCDD	ND	-----	0.0094				
1,2,3,4,6,7,8-HpCDF	ND	-----	0.0130	A	Total 2,3,7,8-TCDD		
1,2,3,4,7,8,9-HpCDF	-----	0.019	0.0140	IA	Equivalence: 0.00020 ng/L		
Total HpCDF	ND	-----	0.0094		(Using ITE Factors)		
1,2,3,4,6,7,8-HpCDD	ND	-----	0.0270	A			
Total HpCDD	ND	-----	0.0094				
OCDF	ND	-----	0.0190				
OCDD	0.20	-----	0.0360	BA			

Conc = Concentration (Totals include 2,3,7,8-substituted isomers)
EMPC = Estimated Maximum Possible Concentration
A = Detection Limit based on signal-to-noise measurement
J = Concentration detected is below the calibration range
B = Less than 10 times higher than method blank level
P = Recovery outside of target range
Nn = Value obtained from additional analysis

LRL = Lower Reporting Limit
I = Interference
E = PCDE Interference
S = Saturated signal
ND = Not Detected
NA = Not Applicable
NC = Not Calculated
* = See Discussion

Report No.....01-1046199

REPORT OF LABORATORY ANALYSIS

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Method 8290 Analysis Results

Client - ROBERT E LEE

Client's Sample ID	01-10563		
Lab Sample ID	102831468		
Filename	U10724A_20		
Injected By	CVS		
Total Amount Extracted	1060 mL	Matrix	WATER
% Moisture	NA	Dilution	NA
Dry Weight Extracted	NA	Collected	06/27/2001
ICAL Date	07/24/2001	Received	06/29/2001
CCal Filename(s)	U10724A_06 & U10724A_21	Extracted	07/10/2001
Method Blank ID	BLANK-1153	Analyzed	07/25/2001 04:41

Native Isomers	Conc ng/L	EMPC ng/L	LRL ng/L		Internal Standards	ng's Added	Percent Recovery
2,3,7,8-TCDF	ND	-----	0.00190	N2	2,3,7,8-TCDF-13C	2.00	79 N2
Total TCDF	ND	-----	0.00190	N2	2,3,7,8-TCDD-13C	2.00	81 N2
					1,2,3,7,8-PeCDF-13C	2.00	97
2,3,7,8-TCDD	ND	-----	0.00190	N2	2,3,4,7,8-PeCDF-13C	2.00	93
Total TCDD	ND	-----	0.00190	N2	1,2,3,7,8-PeCDD-13C	2.00	95
					1,2,3,4,7,8-HxCDF-13C	2.00	97
1,2,3,7,8-PeCDF	ND	-----	0.01000	A	1,2,3,6,7,8-HxCDF-13C	2.00	90
2,3,4,7,8-PeCDF	ND	-----	0.00940		2,3,4,6,7,8-HxCDF-13C	2.00	102
Total PeCDF	ND	-----	0.00940		1,2,3,7,8,9-HxCDF-13C	2.00	97
					1,2,3,4,7,8-HxCDD-13C	2.00	101
1,2,3,7,8-PeCDD	ND	-----	0.01300	A	1,2,3,6,7,8-HxCDD-13C	2.00	92
Total PeCDD	ND	-----	0.00940		1,2,3,4,6,7,8-HpCDF-13C	2.00	96
					1,2,3,4,7,8,9-HpCDF-13C	2.00	102
1,2,3,4,7,8-HxCDF	ND	-----	0.01500	A	1,2,3,4,6,7,8-HpCDD-13C	2.00	105
1,2,3,6,7,8-HxCDF	ND	-----	0.01500	A	OCDD-13C	4.00	106
2,3,4,6,7,8-HxCDF	ND	-----	0.01500	A			
1,2,3,7,8,9-HxCDF	ND	-----	0.01300	A	1,2,3,4-TCDD-13C	2.00	NA
Total HxCDF	ND	-----	0.00940		1,2,3,7,8,9-HxCDD-13C	2.00	NA
1,2,3,4,7,8-HxCDD	ND	-----	0.02500	A	2,3,7,8-TCDD-37Cl4	0.20	84
1,2,3,6,7,8-HxCDD	-----	0.013	0.02300	IA			
1,2,3,7,8,9-HxCDD	ND	-----	0.02700	A			
Total HxCDD	ND	-----	0.00940				
1,2,3,4,6,7,8-HpCDF	-----	0.380	0.01100	EA	Total 2,3,7,8-TCDD		
1,2,3,4,7,8,9-HpCDF	ND	-----	0.02400	A	Equivalence: 0.0077 ng/L		
Total HpCDF	ND	-----	0.00940		(Using ITE Factors)		
1,2,3,4,6,7,8-HpCDD	0.34	-----	0.01800	A			
Total HpCDD	0.52	-----	0.00940				
OCDF	0.32	-----	0.02700	A			
OCDD	3.90	-----	0.04400	A			

Conc = Concentration (Totals include 2,3,7,8-substituted isomers)
EMPC = Estimated Maximum Possible Concentration
A = Detection Limit based on signal-to-noise measurement
J = Concentration detected is below the calibration range
B = Less than 10 times higher than method blank level
P = Recovery outside of target range
Nn = Value obtained from additional analysis

LRL = Lower Reporting Limit
I = Interference
E = PCDE Interference
S = Saturated signal
ND = Not Detected
NA = Not Applicable
NC = Not Calculated
* = See Discussion

Report No.....01-1046199

REPORT OF LABORATORY ANALYSIS

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Method 8290 Laboratory Control Spike Results

Client - ROBERT E LEE

Lab Sample ID	SPIKE-1145	Matrix	WATER
Filename	U10724A_07	Dilution	NA
Total Amount Extracted	1030.6 mL	Extracted	07/10/2001
ICAL Date	07/24/2001	Analyzed	07/24/2001 17:51
CCal Filename(s)	U10724A_06 & U10724A_21	Injected By	CVS
Method Blank ID	BLANK-1153		

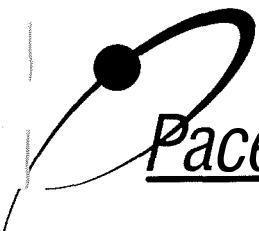
Native Isomers	Qs (ng)	Qm (ng)	% Rec.	Internal Standards	ng's Added	Percent Recovery
2,3,7,8-TCDF	0.20	0.21	103	2,3,7,8-TCDF-13C	2.00	80
				2,3,7,8-TCDD-13C	2.00	83
				1,2,3,7,8-PeCDF-13C	2.00	73
2,3,7,8-TCDD	0.20	0.20	102	2,3,4,7,8-PeCDF-13C	2.00	77
				1,2,3,7,8-PeCDD-13C	2.00	74
				1,2,3,4,7,8-HxCDF-13C	2.00	87
1,2,3,7,8-PeCDF	1.00	1.05	105	1,2,3,6,7,8-HxCDF-13C	2.00	87
2,3,4,7,8-PeCDF	1.00	1.00	100	2,3,4,6,7,8-HxCDF-13C	2.00	96
				1,2,3,7,8,9-HxCDF-13C	2.00	82
				1,2,3,4,7,8-HxCDD-13C	2.00	97
1,2,3,7,8-PeCDD	1.00	1.04	104	1,2,3,6,7,8-HxCDD-13C	2.00	99
				1,2,3,4,6,7,8-HpCDF-13C	2.00	90
				1,2,3,4,7,8,9-HpCDF-13C	2.00	75
1,2,3,4,7,8-HxCDF	1.00	0.99	99	1,2,3,4,6,7,8-HpCDD-13C	2.00	84
1,2,3,6,7,8-HxCDF	1.00	1.05	105	OCDD-13C	4.00	68
2,3,4,6,7,8-HxCDF	1.00	1.01	101			
1,2,3,7,8,9-HxCDF	1.00	1.02	102	1,2,3,4-TCDD-13C	2.00	NA
				1,2,3,7,8,9-HxCDD-13C	2.00	NA
1,2,3,4,7,8-HxCDD	1.00	1.02	102	2,3,7,8-TCDD-37Cl4	0.20	84
1,2,3,6,7,8-HxCDD	1.00	1.00	100			
1,2,3,7,8,9-HxCDD	1.00	0.99	99			
1,2,3,4,6,7,8-HpCDF	1.00	1.02	102			
1,2,3,4,7,8,9-HpCDF	1.00	1.05	105			
1,2,3,4,6,7,8-HpCDD	1.00	1.02	102			
OCDF	2.00	2.05	102			
OCDD	2.00	2.06	103			

Qs = Quantity Spiked
Qm = Quantity Measured
Rec. = Recovery (Expressed as Percent)
P = Recovery outside of target range
X = Background subtracted value
Nn = Value obtained from additional analysis
NA = Not Applicable
* = See Discussion

Report No.....01-1046199

REPORT OF LABORATORY ANALYSIS

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Method 8290 Laboratory Control Spike Results

Client - ROBERT E LEE

Lab Sample ID	SPIKE-DUP-1091	Matrix	WATER
Filename	U10724A_08	Dilution	NA
Total Amount Extracted	1026 mL	Extracted	07/10/2001
ICAL Date	07/24/2001	Analyzed	07/24/2001 18:41
CCal Filename(s)	U10724A_06 & U10724A_21	Injected By	CVS
Method Blank ID	BLANK-1153		

Native Isomers	Qs (ng)	Qm (ng)	% Rec.	Internal Standards	ng's Added	Percent Recovery
2,3,7,8-TCDF	0.20	0.20	98	2,3,7,8-TCDF-13C	2.00	91
				2,3,7,8-TCDD-13C	2.00	87
				1,2,3,7,8-PeCDF-13C	2.00	96
2,3,7,8-TCDD	0.20	0.20	100	2,3,4,7,8-PeCDF-13C	2.00	96
				1,2,3,7,8-PeCDD-13C	2.00	94
				1,2,3,4,7,8-HxCDF-13C	2.00	93
1,2,3,7,8-PeCDF	1.00	0.98	98	1,2,3,6,7,8-HxCDF-13C	2.00	88
2,3,4,7,8-PeCDF	1.00	0.95	95	2,3,4,6,7,8-HxCDF-13C	2.00	101
1,2,3,7,8-PeCDD	1.00	1.00	100	1,2,3,7,8,9-HxCDF-13C	2.00	99
				1,2,3,4,7,8-HxCDD-13C	2.00	96
				1,2,3,6,7,8-HxCDD-13C	2.00	103
1,2,3,4,7,8-HxCDF	1.00	0.96	96	1,2,3,4,6,7,8-HpCDF-13C	2.00	101
1,2,3,6,7,8-HxCDF	1.00	0.97	97	1,2,3,4,7,8,9-HpCDF-13C	2.00	95
2,3,4,6,7,8-HxCDF	1.00	0.93	93	1,2,3,4,6,7,8-HpCDD-13C	2.00	103
1,2,3,7,8,9-HxCDF	1.00	0.96	96	OCDD-13C	4.00	85
1,2,3,4,7,8-HxCDD	1.00	1.03	103	1,2,3,4-TCDD-13C	2.00	NA
				1,2,3,7,8,9-HxCDD-13C	2.00	NA
				2,3,7,8-TCDD-37Cl4	0.20	89
1,2,3,6,7,8-HxCDD	1.00	0.99	99			
1,2,3,7,8,9-HxCDD	1.00	0.98	98			
1,2,3,4,6,7,8-HpCDF	1.00	0.97	97			
				1,2,3,4,7,8,9-HpCDF	1.00	0.99
1,2,3,4,6,7,8-HpCDD	1.00	0.96	96			
OCDF	2.00	1.73	86			
OCDD	2.00	2.04	102			

Qs = Quantity Spiked
 Qm = Quantity Measured
 Rec. = Recovery (Expressed as Percent)
 P = Recovery outside of target range
 X = Background subtracted value
 Nn = Value obtained from additional analysis
 NA = Not Applicable
 * = See Discussion

Report No.....01-1046199

REPORT OF LABORATORY ANALYSIS

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SPIKE RECOVERY RELATIVE PERCENT DIFFERENCE (RPD) RESULTS

Client..... ROBERT E LEE

SPIKE 1 ID..... SPIKE-1145
Spike 1 Filename..... U10724A_07
Spike 2 ID..... SPIKE-DUP-1091
Spike 2 Filename..... U10724A_08

COMPOUND	SPIKE 1 REC,%	SPIKE 2 REC,%	RPD,%
2378-TCDF	103	98	5.0
2378-TCDD	102	100	2.0
12378-PeCDF	105	98	6.9
23478-PeCDF	100	95	5.1
12378-PeCDD	104	100	3.9
123478-HxCDF	99	96	3.1
123678-HxCDF	105	97	7.9
234678-HxCDF	101	93	8.2
123789-HxCDF	102	96	6.1
123478-HxCDD	102	103	1.0
123678-HxCDD	100	99	1.0
123789-HxCDD	99	98	1.0
1234678-HpCDF	102	97	5.0
1234789-HpCDF	105	99	5.9
1234678-HpCDD	102	96	6.1
OCDF	102	86	17.0
OCDD	103	102	1.0

REC = Percent Recovered
RPD = The difference between the two values divided by the average.
NA = Not Applicable

Report No..... 01-1046199

REPORT OF LABORATORY ANALYSIS

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Robert E. Lee & Associates, Inc.

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Green Bay Office 920.336.6338 FAX 920.336.9141
Milwaukee Office 262.569.8893 FAX 262.569.7995

To ensure the proper handling of samples,
please see the back for instructions.

CHAIN OF CUSTODY RECORD

1046199

COC # 85968

Client: Robert E Lee & ASSOC		Project Name: Weisenberger Tie & Lumber		Project Number:		Analyses Required: (Note special detection limits or methods)		Report to: Paul Knuth	
Environmental Program: <input type="checkbox"/> LUST <input type="checkbox"/> SDWA <input type="checkbox"/> WPDES <input type="checkbox"/> RCRA <input type="checkbox"/> OTHER		Requested Turnaround Time <input checked="" type="checkbox"/> Normal (10-15 DAYS) <input type="checkbox"/> Rush		Check Delivery Method <input type="checkbox"/> In Person <input type="checkbox"/> Mail <input type="checkbox"/> Common Courier <input type="checkbox"/> Courier Service <input type="checkbox"/> Other		Date Needed:		Company: Rel	
Sampler:		Sample Type (Matrix) DW = Drinking Water GW = Groundwater WW = Wastewater Soil, Oil, Sludge, Air, Other:		No. Of Containers		Preservation Type (see key below)		Address:	
Date: 6/27/01		Time: 11:00 AM		Received By: Paul Knuth		Date: 6-27-01		Time: 11:00 AM	

Sample Name	Date	Time	Comp	Slab	Filtered	Matrix	No. Of Containers	Preservation Type	REL Sample No.	Remarks
01-10543	6/27/01		A			XW	1	U	102831328	MW-3
10544			P			GW	1	U	336	MW-5
10545			A				1	U	344	MW-6
10546			P				1	U	351	MW-7
10547			A				1	U	369	MW-10
10548			P				1	U	377	DMW-1
10549			A				1	U	385	DMW-2
10550			P				1	U	393	DMW-4
10551			A				1	U	401	DMW-5
10552			P				1	U	419	DMW-6A
10553			A				1	U	427	DMW-7
10554			P				1	U	435	DMW-8

Relinquished By		Date	Time	Received By	Date	Time	Laboratory Receiving Notes Temperature of Contents <u>9</u> °C Custody Seal Intact _____ Sample Condition _____ Sample pH _____
1) _____			A/P	Paul Knuth	6-27-01	11:00 AM	
2) _____			A/P				
3) _____			A/P				

A = AM P = PM

WISCONSIN DNR CERTIFICATION NUMBER 405043870

Preservation Key

N = Nitric Acid O = Sodium Hydroxide
H = Hydrochloric Acid U = Unpreserved
M = Methanol S = Sulfuric Acid



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 Milwaukee Office 262.569.8893 FAX 262.569.7995

To ensure the proper handling of samples,
 please see the back for instructions.

CHAIN OF CUSTODY RECORD

COC # [redacted] 85968

1046199

Client: Robert E Lee & ASSOC

Project Name: _____ Project Number: _____

Weisenberger Tire & Lumber

PO #: _____ BID #: _____

Environmental Program:
 LUST SDWA WPDES RCRA OTHER _____

Requested Turnaround Time: Normal (10-15 DAYS) Rush

Date Needed: _____

Rushes accepted only w/prior notification

Check Delivery Method:
 In Person Mail
 Common Courier Courier Service
 Other _____

Analyses Required:
 (Note special detection limits or methods)

No. Of Containers: _____

Preservation Type (see key below): _____

DIRTY

Report to: Paul Knuth

Company: Rel

Address: _____

Telephone: _____

Fax: _____

Invoice To: _____

Company: _____

Address: _____

Telephone: _____

Fax: _____

Sampler: 3

Sample Type (Matrix)
 DW = Drinking Water
 GW = Groundwater
 WW = Wastewater
 Soil, Oil, Sludge, Air, Other: _____

Sample Name	Date	Time	Comp	Grab	Filter	Notes	REL Sample No.	Remarks
<u>01-10557</u>	<u>6/27/01</u>		<u>A</u>	<u>X</u>	<u>N</u>	<u>GW</u>	<u>10283/4/3</u>	<u>DPZ-1</u>
<u>↓ 10560</u>	<u>↓</u>		<u>A</u>				<u>↓ 450</u>	<u>DPZ-3</u>
<u>↓ 10563</u>	<u>↓</u>		<u>A</u>				<u>↓ 468</u>	<u>DPZ-6</u>
			<u>A</u>					
			<u>P</u>					
			<u>A</u>					
			<u>P</u>					
			<u>A</u>					
			<u>P</u>					
			<u>A</u>					
			<u>P</u>					
			<u>A</u>					
			<u>P</u>					
			<u>A</u>					
			<u>P</u>					
			<u>A</u>					
			<u>P</u>					

Relinquished By	Date	Time	Received By	Date	Time
1) _____	_____	A/P	<u>Ben [signature]</u>	<u>6-27-01</u>	<u>11:00 A/P</u>
2) _____	_____	A/P	_____	_____	_____
3) _____	_____	A/P	_____	_____	_____
Received by Lab _____	_____	_____	_____	_____	_____

Laboratory Receiving Notes

Temperature of Contents 9 °C

Custody Seal Intact _____

Sample Condition _____

Sample pH _____

A = AM P = PM

WISCONSIN DNR CERTIFICATION NUMBER 405043870

Preservation Key
 N = Nitric Acid O = Sodium Hydroxide
 H = Hydrochloric Acid U = Unpreserved
 M = Methanol S = Sulfuric Acid

LABORATORY REPORT
ENVIRONMENTAL HEALTH LABORATORY

- **PRIVATE WELL**

SYNTHETIC ORGANIC ANALYSES

FROM COMMERCIAL LABORATORIES

Section I: To be completed by the Department of Natural Resources

System Name: WEISENBERGER TIE & LUMBER City: _____
PWS ID#: _____ County Code: _____ Route Code: _____
System Well No: _____ Entry Point ID: _____ WI Unique Well No: _____

Sample Point Description: _____

System Type:

Source Code:

Sample Type:

____ (MC) Municipal Community _____ W Well _____ D (SDWA) Compliance Sample
____ (OC) OTM Community _____ E Entry Point _____ C (SDWA) Confirmation _____ - _____ - _____
____ (NN) Nontransient Noncommunity _____ D Distribution (Initial Sample Date)
____ (TN) Transient Noncommunity _____ W Raw Water Sample
____ I Investigation Sample

Collect sample by: _____ - _____ - _____ Return results to DNR by: _____ - _____ - _____

Section II: To be completed by SAMPLER

Sample Collection Date: 06 - 27 - 2001 Sample Collection Time: _____ : _____

Sample Point Address: _____

Sample Point Descrip: _____

First Initial and
Last Name of Sampler: _____ - _____

Section III: To be completed by LABORATORY OFFICIAL. Report analytical results on back.

Laboratory ID Number: 999766900 Laboratory Name: Environmental Health Laboratories

Date Sample Received: 06 - 29 - 2001 Time Sample Received: 09 : 15 Laboratory Sample ID: 624194

Signature of Receiving Lab Official: *Computer for Mike Love* Date Reported: 08 - 07 - 2001

Condition of Sample Upon Receipt: 8° C

Section IV: To be completed by WATER SUPPLY SYSTEM OFFICIAL after analysis has been done.

I certify that I have personally examined and am familiar with the information submitted on this document and all attachments and that, based on my inquiry of those individuals immediately responsible for obtaining the information, I believe that the information is true and accurate, and complete. I also certify that the values being submitted are the actual values found in the sample; no values have been modified or changed in any manner.

Signature: _____ Title: _____

Date Signed: _____

This page to be completed by WATER SUPPLY SYSTEM OFFICIAL or laboratory performing analysis.

Storet Code	Parameter	SDWA Method	MDL *	Results	MCL	Units
46317	Alachlor (Lasso)				2	ug/L
39053	Aldicarb				3	ug/L
82587	Aldicarb Sulfone				2	ug/L
82576	Aldicarb Sulfoxide				4	ug/L
34680	Aldrin				---	ug/L
39033	Atrazine				3	ug/L
34247	Benzo(a)pyrene				0.2	ug/L
77860	Butachlor				---	ug/L
77700	Carbaryl				---	ug/L
81405	Carbofuran				40	ug/L
39348	Chlordane alpha				---	ug/L
39810	Chlordane gamma				---	ug/L
39350	Chlordane				2	ug/L
39730	X 2,4-D	515.1	0.1	< 0.1	70	ug/L
38432	X Dalapon	515.1	1.0	< 1.0	200	ug/L
46373	Deethylatrazine				---	ug/L
46374	Deisopropylatrazine				---	ug/L
99075	Diaminoatrazine				---	ug/L
38760	1,2-Dibromo-3-chloropropane (DBCP)				0.2	ug/L
82052	X Dicamba	515.1	0.1	< 0.1	---	ug/L
39380	Dieldrin				---	ug/L
77903	Di(2-ethylhexyl)adipate				400	ug/L
46312	Di(2-ethylhexyl)phthalate				6	ug/L
81287	X Dinoseb	515.1	0.1	< 0.1	7	ug/L
78885	Diquat				20	ug/L
38926	Endothall				100	ug/L
39390	Endrin				2.0	ug/L
46396	Ethylene dibromide (EDB)				0.05	ug/L
39941	Glyphosate (Round-up)				700	ug/L
39410	Heptachlor				0.4	ug/L
39420	Heptachlor epoxide				0.2	ug/L
34688	Hexachlorobenzene				1	ug/L
34386	Hexachlorocyclopentadiene				50	ug/L
82584	3-Hydroxycarbofuran				---	ug/L
39340	BHC gamma (Lindane)				0.2	ug/L
39480	Methoxychlor				40	ug/L
39051	Methomyl				---	ug/L
39356	Dual (Metolachlor)				---	ug/L
81408	Metribuzin (Sencor)				---	ug/L
38865	Oxamyl (Vydate)				200	ug/L
39515	PCB Total				0.5	ug/L
39032	X Pentachlorophenol	515.1	0.04	0.05	1	ug/L
39720	X Picloram (Tordon)	515.1	0.1	< 0.1	500	ug/L
30295	Propachlor				---	ug/L
39760	X 2,4,5-TP (Silvex)	515.1	0.1	< 0.1	50	ug/L
39055	Simazine				4	ug/L
34675	2,3,7,8-TCDD (Dioxin)				0.00003	ug/L
39400	Toxaphene				3	ug/L

* EHL has demonstrated it can achieve these report limits in reagent water, but can not document them in all sample matrices.

Site:

Report #: 624194

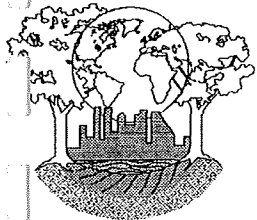
REPORT SUMMARY

1,2,4-Trichlorophenol was detected in the sample submitted for analysis at a concentration of 0.05 ug/L, which is less than the current MCL of 1 ug/L. None of the other analytes included in the detailed parameter list were detected in the sample submitted for analysis.

Note: The sample submitted for analysis was received at 8° C. The client was notified of the situation, and analysis was authorized by Paul Knuth of Robert E. Lee & Associates, Inc.

Note: Sample container was provided by the client. REL Sample Name 01-10567

Analysis Date: 08/01/2001



Laboratory Name: Environmental Health Laboratories Laboratory ID Number: 999766900

Note: This report may not be reproduced, except in full, without written approval from Environmental Health Laboratories(EHL). EHL is a National Environmental Laboratory Accreditation Program (NELAP) accredited laboratory. These results may not meet the NELAP standard but satisfy the requirements for this project.

We appreciate the opportunity to provide you with this analysis. If you have any questions concerning this report, please do not hesitate to call us at (219) 233-4777.

Reviewed By: Steve Dungey

Date: 8/7/01

Analized By: [Signature]

Date: 8-8-01

