

10/29/96

246100800ERR-LUST
03-46-003301 h

**THIRD QUARTER 1996
GROUND-WATER QUALITY
MONITORING**

**FORMER CEDARBURG LIGHT AND
WATER COMMISSION POWER PLANT
W61 N617 MEQUON AVENUE
CEDARBURG, WISCONSIN**

**(WDNR FID #246100800 ERR-LUST)
(PECFA CLAIM #53012-2017-17)**

October 25, 1996

October 25, 1996
(CLW131246.700)

Mr. Dale Lythjohan
Cedarburg Light and Water Commission
N30 W5926 Lincoln Boulevard
Post Office Box 767
Cedarburg, Wisconsin 53012

RE: Third Quarter 1996 Ground-Water Quality Monitoring, Former Cedarburg Light and Water Commission Power Plant, W61 N617 Mequon Avenue, Cedarburg, Wisconsin (WDNR FID #246100800ERR-LUST) (PECFA Claim #53012-2017-17)

Dear Mr. Lythjohan:

Northern Environmental Technologies, Incorporated (Northern Environmental) completed the third round of the Wisconsin Department of Natural Resources (WDNR)-approved four rounds of quarterly ground-water quality monitoring at the former Cedarburg Light and Water Commission Power Plant, W61 N617 Mequon Avenue, Cedarburg, Wisconsin (the Property). The Property is in the Southeast Quarter of the Northeast Quarter of Section 27, Township 10 North, Range 21 East in Ozaukee County, Wisconsin (latitude 43 degrees, 18 minutes, 22 seconds north, longitude 87 degrees, 59 minutes, 40 seconds west) (Figure 1) (USGS, 1976).

BACKGROUND INFORMATION

The Property was formerly an electrical generating plant owned and operated by the Cedarburg Light and Water Commission. Two 20,000-gallon capacity diesel fuel underground storage tanks (USTs) were reportedly cleaned and abandoned in place at the Property during April 1986. One 1000-gallon gasoline/diesel UST was also cleaned, removed, and disposed at that time. A closure assessment was not required when the USTs were decommissioned.

During 1993, the Cedarburg Light and Water Commission retained Northern Environmental to drill and sample boreholes on the Property as part of an environmental assessment (Northern Environmental, 1993). Diesel range organics (DRO) and gasoline range organics (GRO) were detected in soil samples. Northern Environmental completed a site investigation during April 1994. Contaminated ground water was discovered. A report was prepared that described the investigation and presented the results (Northern Environmental, 1994). The WDNR requested that an additional monitoring well be installed south of the power plant. The well (MW500) was drilled and installed during December 1994. No DRO or petroleum volatile organic

compounds (PVOCS) were detected in the soil sample collected from the monitoring well borehole. No volatile organic compounds (VOCs) or DRO were detected in water from the new well.

In an unrelated remedial action, Mercury Marine, Incorporated removed polychlorinated biphenyl-contaminated sediments from Ruck Pond during 1994. DRO-contaminated soil was discovered in stream bank excavations abutting the Property. The WDNR and Cedarburg Light and Water Commission were notified. Northern Environmental collected soil samples from the excavations to assess the extent of contamination. The soil samples were laboratory analyzed for DRO and PVOCS. High concentrations of DRO are present in soil beneath the cooling towers on the bank of Ruck Pond.

Additional ground-water quality monitoring was performed during January and June 1995. Ground-water samples from MW200 contained trichloroethene and tetrachloroethene above the Chapter NR 140, Wisconsin Administrative Code (NR 140, Wis. Adm. Code) enforcement standards (ES). Benzene concentrations above the NR 140, Wis. Adm. Code preventive action limit (PAL) were detected in samples from MW300. A report was prepared describing the excavation sampling and additional ground-water monitoring (Northern Environmental, 1995).

The WDNR approved long-term ground-water monitoring as an appropriate remedial response (Vance, 1996). The WDNR requested that the monitoring wells be sampled quarterly for one year with sampling frequency reduced to once per year thereafter depending on results (Vance, 1996). This letter describes the third round of quarterly ground-water quality sampling.

METHODS OF INVESTIGATION

Ground-water samples were collected from the four monitoring wells on September 13, 1996 to monitor ground-water quality at the Property. Before purging and sampling the monitoring wells, Northern Environmental personnel measured the depth to water in each well to evaluate ground-water flow direction. The depth to water measurements were converted to elevations relative to a site datum. The monitoring wells were purged before sampling in accordance with WDNR requirements (NR 141, Wis. Adm. Code).

Ground-water samples were collected by gently lowering new bottom-filling disposable polyethylene bailers into the wells until the bailer was completely submerged. Water samples were transferred from the bailers into appropriate sample containers using new bottom-emptying devices. The samples were preserved with hydrochloric acid, labeled, and chilled until delivery to U.S. Analytical Laboratory (Combined Locks, Wisconsin) (WDNR Certification #445027660) for analysis. The ground-water samples were laboratory analyzed for DRO using the WDNR Modified Method and VOCs using Environmental Protection Agency Method 8021.

Quality assurance/quality control (QA/QC) samples were also collected in accordance with WDNR guidelines. QA/QC samples consisted of one trip blank, one field blank, and one duplicate sample. The trip blank (labeled "Trip Blank") was obtained from the laboratory and accompanied the investigative samples throughout the chain-of-custody. The field blank (sample labeled FB-1246) was prepared with laboratory-grade deionized water using the same

techniques and equipment as the investigative samples. The duplicate sample (sample labeled Dup-1246) was collected from MW200. QA/QC samples were analyzed for VOCs using the before mentioned method. Ground-water sample laboratory reports and the associated chain-of-custody record are provided in Attachment A.

SUMMARY OF FINDINGS

DRO and VOCs were not detected in monitoring well MW500. DRO was present in the ground-water sample from MW200. Tetrachloroethene concentrations exceeded the ES, and trichloroethene and benzene exceeded the PAL in MW200. Tetrachloroethene exceeded the PAL in MW300 and MW400. All other VOCs were either not detected or were below their respective PALs. Ground-water sample laboratory analysis results are summarized in Table 2.

June 10, 1996 water table conditions are depicted in Figure 2. Ground-water flow was predominantly eastward across the Property under an approximate hydraulic gradient of 0.01 foot per foot. Water table elevation data is summarized in Table 1.

CONCLUSIONS AND RECOMMENDATIONS

Benzene, tetrachloroethene, and trichloroethene were present in monitoring wells MW200 and MW300. Tetrachloroethene in MW200 was the only compound exceeding the ES. Ground-water quality during this round of sampling is generally consistent with previous results. The final round of the WDNR-approved water quality monitoring will be performed during December 1996. The ground-water quality data collected over the past year will then be evaluated to determine a new ground-water monitoring sampling schedule.

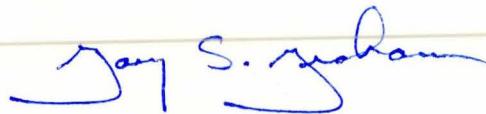
The results of this study are based upon professional interpretation of the information available to Northern Environmental given site conditions and the time and budget constraints of this project. Northern Environmental does not warrant that this report represents an exhaustive study of all possible impacts at the study area. The items investigated as part of this investigation do represent the most likely sources of environmental impacts associated with the described UST systems, and are consequently believed to adequately address WDNR requirements and the needs of the client at the present time.

We trust this information meets your needs. Please feel free to contact us if you have any questions or comments.

Sincerely,
**Northern Environmental
Technologies, Incorporated**



Christopher C. Hatfield
Hydrogeologist



Gary S. Graham
Senior Project Manager

CCH/kmr

Enclosures

cc: Ms. Kaye Vance (Cook and Franke, S.C.)
Mr. John Feeney (Wisconsin Department of Natural Resources)

REFERENCES

Northern Environmental Technologies, Incorporated, "Phase I Environmental Site Assessment, Cedarburg Light and Water Commission Former Power Plant, W61 N617 Mequon Avenue, Cedarburg, Wisconsin," February 4, 1993.

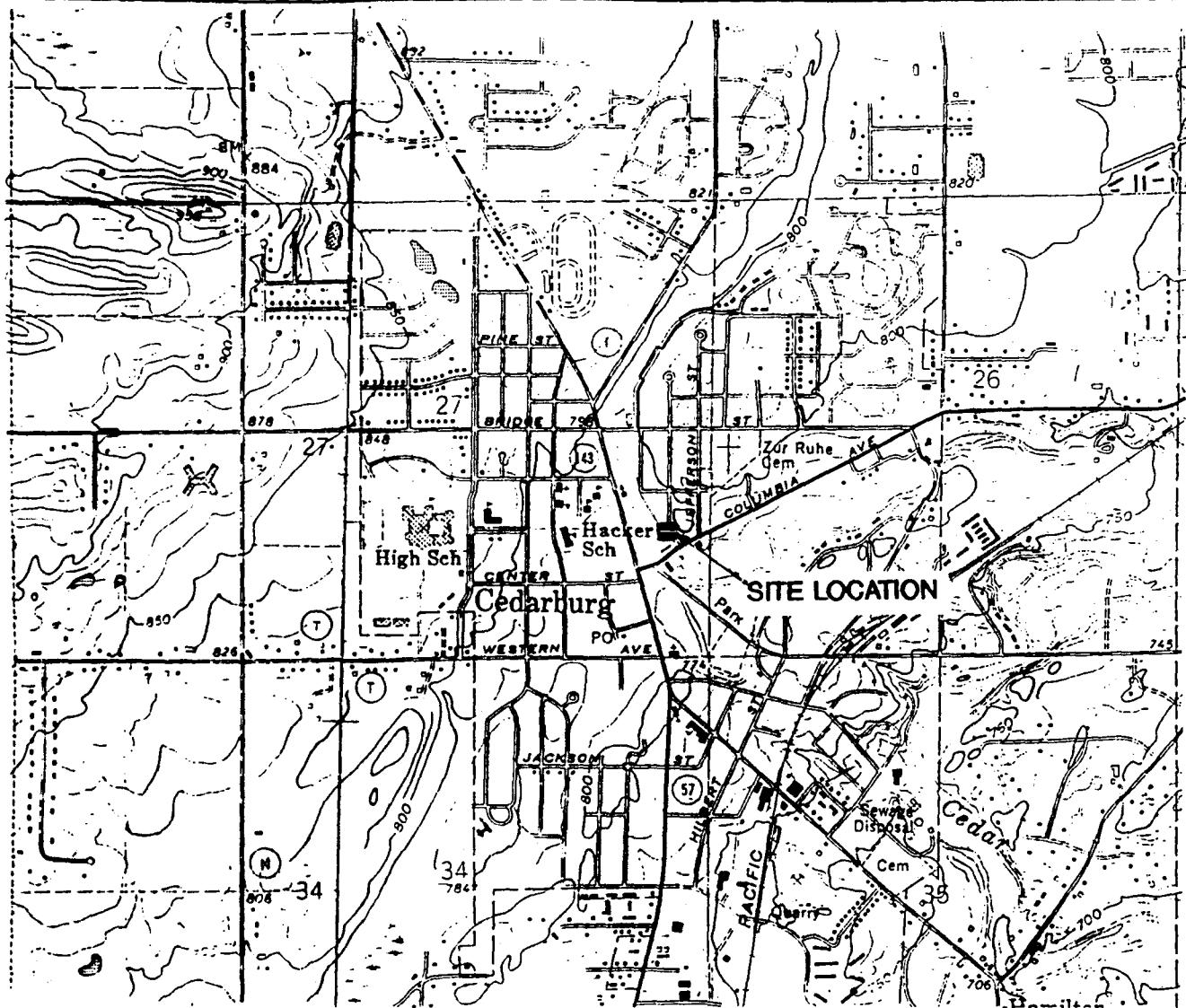
Northern Environmental Technologies, Incorporated, "Site Investigation Results, Former Cedarburg Light and Water Commission Power Plant, W61 N617 Mequon Avenue, Cedarburg, Wisconsin," October 19, 1995.

Northern Environmental Technologies, Incorporated, "Site Investigation Results, Former Power Plant, W61 N617 Mequon Avenue, Cedarburg, Wisconsin," April 15, 1994.

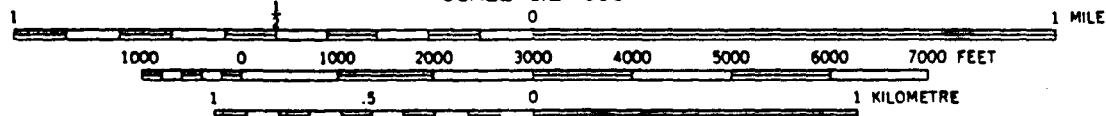
United States Geological Survey, *Cedarburg, Wisconsin, 7.5 Minute Quadrangle Topographic Map*, 1959, Photorevised 1971 and 1976.

Vance, Kaye (Cedarburg City Attorney), letter to John Feeney (Wisconsin Department of Natural Resources), February 13, 1996.

Wisconsin Department of Natural Resources, "Ground-Water Monitoring Well Requirements," *Wisconsin Administrative Code*, Chapter NR 141, June 1991.



SCALE 1:24 000



CONTOUR INTERVAL 10 FEET
NATIONAL GEODETIC VERTICAL DATUM OF 1929



BASE MAP SOURCE: USGS CEDARBURG AND FIVE CORNERS, WI 7.5 MIN QUADRANGLE

| | | | |
|---|--------------------|---------------|--|
| REV | PROJECT: CLW131246 | DATE 10/25/96 | CEDARBURG LIGHT & WATER COMMISSION CEDARBURG, WISCONSIN |
| <p>THIS DRAWING AND ALL INFORMATION CONTAINED THEREON IS THE PROPERTY OF NORTHERN ENVIRONMENTAL INCORPORATED AND SHALL NOT BE COPIED OR USED EXCEPT FOR THE PURPOSE FOR WHICH IT IS EXPRESSLY FURNISHED. THE DRAWING AND ANY COPIES THEREOF SHALL BE RETURNED TO THE OWNER ON DEMAND.</p> | | | |
| <p>▲ Northern Environmental Hydrologists • Engineers • Geologists</p> | | | SITE LOCATION AND LOCAL TOPOGRAPHY |

FIGURE 1

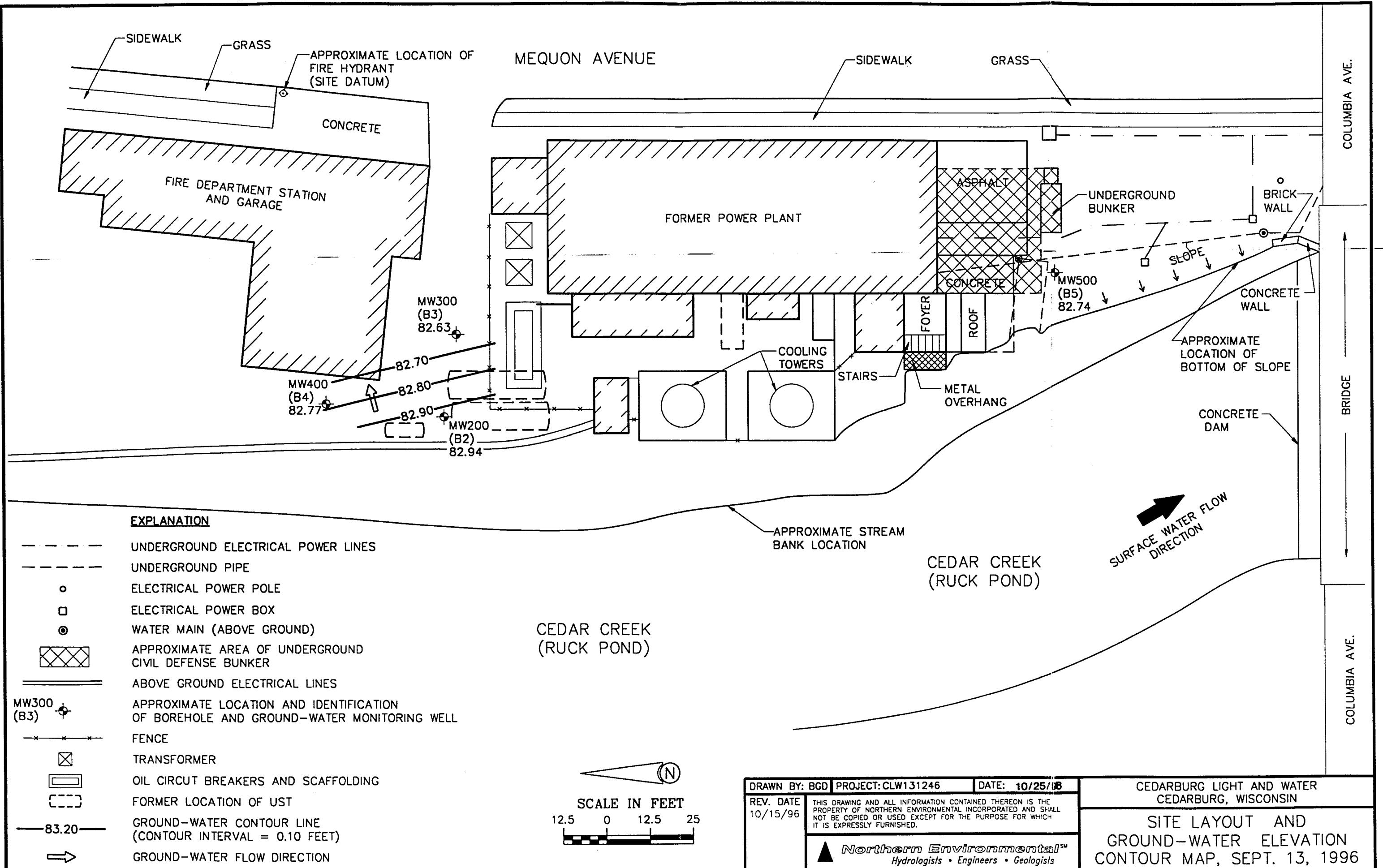


Table 1 Ground-Water Elevation Data, Former Power Plant, Cedarburg, Wisconsin

| Well ID | Elevation Ground Surface (feet) | Elevation of Reference Point* (feet) | Date | Depth to Water Below Reference Point* (feet) | Water Table Elevation (feet) |
|---------|---------------------------------------|--|----------|--|------------------------------------|
| MW200 | 96.46 | 95.94 | 10/18/93 | 12.71 | 83.23 |
| | | | 10/25/93 | 12.78 | 83.16 |
| | | | 10/28/93 | 12.94 | 83.00 |
| | | | 01/07/94 | 13.30 | 82.64 |
| | | | 02/14/94 | 14.21 | 81.73 |
| | | | 12/28/94 | 13.02 | 82.92 |
| | | | 01/18/95 | 12.90 | 83.04 |
| | | | 06/08/95 | 12.53 | 83.41 |
| | | | 03/21/96 | 12.81 | 83.13 |
| | | | 06/10/96 | 11.49 | 84.45 |
| | | | 09/13/96 | 13.00 | 82.94 |
| MW300 | 97.22 | 96.54 | 10/18/93 | 14.02 | 82.52 |
| | | | 10/25/93 | 14.01 | 82.53 |
| | | | 10/28/93 | 13.98 | 82.56 |
| | | | 01/07/94 | 14.41 | 82.13 |
| | | | 02/14/94 | 15.16 | 81.38 |
| | | | 12/28/94 | 14.01 | 82.53 |
| | | | 01/18/95 | 12.91 | 83.63 |
| | | | 06/08/95 | 13.42 | 83.12 |
| | | | 03/21/96 | 13.76 | 82.78 |
| | | | 06/10/96 | 12.31 | 84.23 |
| | | | 09/13/96 | 13.91 | 82.63 |
| MW400 | 95.56 | 95.28 | 10/18/93 | 12.60 | 82.68 |
| | | | 10/25/93 | 12.58 | 82.70 |
| | | | 10/28/93 | 12.55 | 82.73 |
| | | | 01/07/94 | 12.87 | 82.41 |
| | | | 02/14/94 | 13.62 | 81.66 |
| | | | 12/28/94 | 12.50 | 82.78 |
| | | | 01/18/95 | 12.38 | 82.90 |
| | | | 06/08/95 | 12.03 | 83.25 |
| | | | 03/21/96 | 12.28 | 83.00 |
| | | | 06/10/96 | 10.96 | 84.32 |
| | | | 09/13/96 | 12.51 | 82.77 |
| MW500 | 95.53** | 95.56 | 12/28/94 | 12.54 | 83.02 |
| | | | 01/03/95 | 12.42 | 83.14 |
| | | | 01/18/95 | 12.42 | 83.14 |
| | | | 06/08/95 | 12.62 | 82.94 |
| | | | 03/21/96 | 12.62 | 82.94 |
| | | | 06/10/96 | 11.42 | 84.14 |
| | | | 09/13/96 | 12.82 | 82.74 |

NOTE: Elevations are referenced to site datum

* = Reference point is the top of the monitoring well casing

** = Elevation of top of protective metal casing

Table 2 Ground-Water Analysis Results, Former Power Plant, Cedarburg, Wisconsin

| Well I.D. | Date | Concentrations of Detected Analytes (µg/l) | | | | | | | | | | | | | | | | | | |
|-------------|-----------|--|-------|---------|---------------|---------|---------------|-------|-----------------|---------------|---------------------|-------------------------|-------------|---------------------|------------------|--------------------------|--------------------------|----------------------|------------------------|------|
| | | DRO | GRO | Benzene | Ethyl-benzene | Toluene | Total Xylenes | MTBE | n-butyl-benzene | chloro-ethane | 1,1-Dichloro-ethane | cis-1,2-Dichloro-ethene | Naphthalene | Tetra-chloro-ethene | Trichloro-ethene | 1,2,4-Tri-methyl-benzene | 1,3,5-Tri-methyl-benzene | 1,2-Dichloro-benzene | 1,1,1-Trichloro-ethane | Lead |
| MW200 | 10/28/93 | 720 | 110 | <0.6 | <1.0 | 35 | 5.6 | <1.0 | 6.1 | 23 | 7.4 | 3.5 | 5.7 | 5.4 | 7.6 | 5.7 | 3.2 | <1.0 | <0.2 | 17 |
| | 01/13/94 | <5.0 | <10.0 | <0.6 | <1.0 | 2.4 | 1.8 | <1.0 | 3.0 | 26 | 3.6 | 1.2 | 7.1 | 1.4 | 1.6 | 2.1 | 1.5 | 1.6 | <0.2 | 22 |
| | 01/18/95 | 2000 | 28 | <2.0 | <1.0 | <1.0 | <2.0 | <1.0 | <2.0 | 2.2 | 4.9 | 22 | 0.44 | 19 | 29 | <2.0 | 0.19 | 4.9 | 4 | |
| | *01/18/95 | NA | NA | 0.28 | <1.0 | <1.0 | <2.5 | <1.0 | <2.0 | 3.1 | 4.1 | 22 | <2.0 | 20 | 30 | <1.0 | <1.0 | 5.0 | NA | |
| | 06/08/95 | 810 | NA | <.26 | <.32 | <.69 | <1.23 | .46 | <.45 | 9.4 | 6.6 | 8.4 | <.41 | 42 | 17.5 | <.57 | .33 | 2.9 | <1 | |
| | 03/21/96 | 510 | NA | 0.28 | <0.32 | <0.69 | <1.23 | 0.29 | <0.45 | 6.1 | 4.1 | 5.5 | <0.41 | 11 | 5.8 | <0.57 | 0.69 | 0.65 | NA | |
| | *03/21/96 | NA | NA | 0.28 | <0.32 | <0.69 | <1.23 | 0.34 | 0.69 | 6.6 | 4.0 | 4.8 | 0.68 | 9.5 | 5.1 | <0.57 | 0.76 | <0.63 | NA | |
| | 06/10/96 | 270 | NA | 0.27 | <0.32 | <0.69 | <1.23 | <0.22 | <0.45 | 6.2 | 5.9 | 6.9 | <0.41 | 56 | 14 | <0.57 | 0.43 | 2.8 | NA | |
| | *06/10/96 | NA | NA | 0.29 | <0.32 | <0.69 | <1.23 | <0.22 | <0.45 | 6.8 | 5.9 | 6.8 | <0.41 | 51 | 14 | <0.57 | 0.51 | 2.8 | NA | |
| | 09/13/96 | 350 | NA | 0.48 | <0.32 | <0.69 | <1.23 | <0.22 | <0.45 | 2.6 | 4.7 | 3.9 | <0.41 | 15 | 4.7 | <0.57 | 0.93 | 0.97 | NA | |
| | *09/13/96 | NA | NA | 0.52 | <0.32 | <0.69 | <1.23 | <0.22 | <0.45 | 2.6 | 4.9 | 4.1 | <0.41 | 15 | 4.8 | <0.57 | 1.0 | 1.0 | NA | |
| MW300 | 10/28/93 | <100 | <100 | 1.2 | | 1.5 | <2.5 | | <2.0 | 3.3 | 5.0 | 3.4 | <2.0 | 3.9 | <1.0 | <1.0 | <1.0 | <1.0 | 2 | |
| | 01/13/94 | <5.0 | <10.0 | 1.3 | <1.0 | <1.0 | <2.5 | <1.0 | <2.0 | <1.0 | <1.0 | <1.0 | <2.0 | <1.0 | <1.0 | <1.0 | <1.0 | <0.2 | <1.0 | |
| | 01/18/95 | 150 | <11.0 | 0.80 | <1.0 | <1.0 | <2.5 | <1.0 | <2.0 | 2.3 | 1.1 | 0.90 | <2.0 | <1.0 | <1.0 | <1.0 | <1.0 | <1.0 | 1.0 | |
| | 06/08/95 | <100 | NA | 0.36 | <.32 | <.69 | <1.23 | <.22 | <.45 | 0.93 | 0.9 | 0.67 | <.41 | 1.82 | 0.33 | <.57 | <.57 | <.63 | 1.0 | |
| | 03/21/96 | 400 | NA | 1.1 | <0.32 | <0.69 | 0.62 | <0.22 | <0.45 | 3.9 | <0.37 | 0.32 | 4.9 | 1.5 | <0.18 | <0.57 | 0.31 | <0.63 | NA | |
| | 06/10/96 | <100 | NA | 0.41 | <0.32 | <0.69 | <1.23 | <0.22 | <0.45 | 1.3 | 0.75 | 0.75 | 1.7 | 2.1 | 0.45 | <0.57 | 0.12 | <0.63 | NA | |
| | 09/13/96 | <100 | NA | 0.34 | <0.32 | <0.69 | <1.23 | <0.22 | <0.45 | 1.5 | 0.63 | 0.59 | 0.56 | 3.2 | 0.49 | <0.57 | 0.14 | <0.63 | NA | |
| MW400 | 10/28/93 | <100 | <100 | <0.6 | <1.0 | <1.0 | <2.5 | <1.0 | <2.0 | <1.0 | <1.0 | <1.0 | <2.0 | <1.0 | <1.0 | <1.0 | <1.0 | <0.2 | <1.0 | |
| | 01/13/94 | <5.0 | <10.0 | <0.6 | <1.0 | <1.0 | <2.5 | <1.0 | <2.0 | <1.0 | <1.0 | <1.0 | <2.0 | <1.0 | <1.0 | <1.0 | <1.0 | <0.2 | <1.0 | |
| | 01/18/95 | 120 | <11.0 | <0.6 | <1.0 | <1.0 | <2.5 | <1.0 | <2.0 | <1.0 | <1.0 | <1.0 | <2.0 | <1.0 | <1.0 | <1.0 | <1.0 | <0.2 | 1.0 | |
| | 06/08/95 | <100 | NA | <.26 | <.32 | <.69 | <1.23 | .33 | <.45 | <.5 | <.27 | <.29 | <.41 | <.56 | <.18 | <.57 | <.57 | <.63 | 2 | |
| | 03/21/96 | <100 | NA | <0.26 | <0.32 | <0.69 | <1.23 | <0.22 | <0.45 | <0.5 | <0.37 | <0.29 | <0.41 | <0.56 | <0.18 | <0.57 | <0.57 | <0.63 | NA | |
| | 06/10/96 | <100 | NA | <0.28 | <0.32 | <0.69 | <1.23 | <0.22 | <0.45 | <0.5 | <0.27 | <0.29 | <0.41 | <0.56 | <0.18 | <0.57 | <0.57 | <0.63 | NA | |
| | 09/13/96 | <100 | NA | <0.26 | <0.32 | <0.69 | <1.23 | <0.22 | <0.45 | <0.5 | <0.27 | <0.29 | <0.41 | 0.68 | <0.18 | <0.57 | <0.57 | <0.63 | NA | |
| MW500 | 01/18/95 | <100 | <11 | <0.6 | <1.0 | <1.0 | <2.5 | <1.0 | <2.0 | <1.0 | <1.0 | <1.0 | <2.0 | <1.0 | <1.0 | <1.0 | <1.0 | <0.2 | <1.0 | |
| | 06/08/95 | <100 | NA | <.26 | <.32 | <.69 | <1.23 | <.22 | <.45 | <.5 | <.27 | <.29 | <.41 | <.56 | <.18 | <.57 | <.57 | <.63 | 2 | |
| | 03/21/96 | <100 | NA | <0.26 | <0.32 | <0.69 | <1.23 | <0.22 | <0.45 | <0.5 | <0.37 | <0.29 | <0.41 | <0.56 | <0.18 | <0.57 | <0.57 | <0.63 | NA | |
| | 06/10/96 | <100 | NA | <0.28 | <0.32 | <0.69 | <1.23 | <0.22 | <0.45 | <0.5 | <0.27 | <0.29 | <0.41 | <0.56 | <0.18 | <0.57 | <0.57 | <0.63 | NA | |
| | 09/13/96 | <100 | NA | <0.28 | <0.32 | <0.69 | <1.23 | <0.22 | <0.45 | <0.5 | <0.27 | <0.29 | <0.41 | <0.56 | <0.18 | <0.57 | <0.57 | <0.63 | NA | |
| Field Blank | 01/18/95 | NA | NA | <0.6 | <1.0 | <1.0 | <2.5 | <1.0 | <2.0 | <1.0 | <1.0 | <1.0 | <2.0 | <1.0 | <1.0 | <1.0 | <1.0 | <0.2 | NA | |
| | 06/08/95 | NA | NA | <.26 | <.32 | <.69 | <1.23 | <.22 | <.45 | <.5 | <.27 | <.29 | <.41 | <.56 | <.18 | <.57 | <.57 | <.63 | NA | |
| | 03/21/96 | NA | NA | <0.26 | <0.32 | <0.69 | <1.23 | <0.22 | <0.45 | <0.5 | <0.37 | <0.29 | <0.41 | <0.56 | <0.18 | <0.57 | <0.57 | <0.63 | NA | |
| | 06/10/96 | <100 | NA | <0.28 | <0.32 | <0.69 | <1.23 | <0.22 | <0.45 | <0.5 | <0.27 | <0.29 | <0.41 | <0.56 | <0.18 | <0.57 | <0.57 | <0.63 | NA | |
| Trip Blank | 06/10/96 | <100 | NA | <0.28 | <0.32 | <0.69 | <1.23 | <0.22 | <0.45 | <0.5 | <0.27 | <0.29</td | | | | | | | | |

ATTACHMENT A

LABORATORY REPORTS AND
ASSOCIATED CHAIN-OF-CUSTODY RECORD

**Analytical Laboratory**

1090 Kennedy Ave. Kimberly, WI 54136
414-735-8295

WI DNR Certified Lab #445027660

BRIAN KRANZ
NORTHERN ENVIRONMENTAL
1214 W VENTURE COURT
MEQUON WI 53092

Project #: CLW131246
Project : Cedarburg
Sample ID: MW200
Lab Code: 5014513A
Sample Type: Water
Sample Date: 13-Sep-96

Report Date: 02-Oct-96

| Test | Result | MDL | PQL | Unit | pH | Date Ext/Digested | Date Analyzed: | Analyzed By: | QC Code |
|-----------------------------|--------|-----|-----|------|-----|-------------------|----------------|--------------|---------|
| MODIFIED DRO WDNR SEP 95 | 350 | 30 | 96 | UG/L | 2.5 | 19-Sep-96 | 19-Sep-96 | C. Rotar | 1 |

MDL = Method Detection Limit

PQL = Practical Quantitation Limit

QC SUMMARY

CODE:

1 pH adjusted below two.

Authorized Signature



Analytical Laboratory

1090 Kennedy Ave. Kimberly, WI 54136
414-735-8295

WI DNR Certified Lab #445027660

Method 8021 Volatile Organic Compounds

BRIAN KRANZ
NORTHERN ENVIRONMENTAL
1214 W VENTURE COURT
MEQUON WI 53092

Report Date: 03-Oct-96
Analyzed By: G.Shah

Project #: CLW131246
Project : Cedarburg
Sample ID: MW200
Lab Code: 5014513A
Sample Type: Water
Sample Date: 13-Sep-96
Date Analyzed: 25-Sep-96

| ANALYTE | RESULT | MDL UG/L | PQL UG/L |
|-----------------------------|--------|-------------|-------------|
| Benzene | 0.48 | 0.082 | 0.26 |
| Bromobenzene | < 0.24 | 0.075 | 0.24 |
| Bromodichloromethane | < 0.11 | 0.035 | 0.11 |
| n-Butylbenzene | < 0.45 | 0.14 | 0.45 |
| sec-Butylbenzene | 0.6 | 0.15 | 0.49 |
| tert-Butylbenzene | < 0.4 | 0.12 | 0.4 |
| Carbon Tetrachloride | < 0.5 | 0.16 | 0.5 |
| Chlorobenzene | < 0.27 | 0.086 | 0.27 |
| Chloroethane | 2.6 | 0.39 | 1.3 |
| Chloroform | < 0.22 | 0.07 | 0.22 |
| Chloromethane | < 1 | 0.88 | 3.1 |
| 2-Chlorotoluene | < 0.65 | 0.21 | 0.65 |
| 4-Chlorotoluene | < 0.19 | 0.06 | 0.19 |
| 1,2-Dibromo-3-Chloropropane | < 1 | 0.83 | 2.7 |
| Dibromochloromethane | < 0.09 | 0.028 | 0.09 |
| 1,2-Dichlorobenzene | 0.93 | 0.035 | 0.11 |
| 1,3-Dichlorobenzene | < 0.83 | 0.23 | 0.83 |
| 1,4-Dichlorobenzene | < 0.13 | 0.039 | 0.13 |
| Dichlorodifluoromethane | < 5.4 | 1.7 | 5.4 |
| 1,1-Dichloroethene | < 0.37 | 0.12 | 0.37 |
| 1,2-Dichloroethane | < 0.86 | 0.27 | 0.86 |
| 1,1-Dichloroethane | 4.7 | 0.084 | 0.27 |
| cis 1,2-Dichloroethene | 3.9 | 0.092 | 0.29 |
| trans-1,2-dichloroethene | 0.28 | 0.072 | 0.23 |
| 1,2-Dichloropropane | < 0.15 | 0.046 | 0.15 |
| 1,3-DCP, Tetrachloroethene | < 0.56 | 0.17 | 0.56 |

Fluorobenzene Surrogate 107 % Rec.
1,4-Dichlorobutane Surrogate 84 % Rec.
Sample pH 1.6

| ANALYTE | RESULT | MDL UG/L | PQL UG/L |
|---------------------------|--------|-------------|-------------|
| 2,2-Dichloropropane | < 1 | 0.63 | 2.2 |
| Di-isopropyl Ether | < 0.38 | 0.12 | 0.38 |
| Ethylbenzene | < 0.32 | 0.1 | 0.32 |
| EDB (1,2-Dibromoethane) | 0.16 | 0.025 | 0.08 |
| Hexachlorobutadiene | < 0.35 | 0.11 | 0.35 |
| Isopropylbenzene | < 0.36 | 0.11 | 0.36 |
| p-Isopropyltoluene | < 0.46 | 0.15 | 0.46 |
| Methylene Chloride | < 4 | 0.29 | 0.91 |
| MTBE | < 0.22 | 0.069 | 0.22 |
| Naphthalene | < 0.41 | 0.13 | 0.41 |
| n-Propylbenzene | < 0.41 | 0.13 | 0.41 |
| 1,1,2,2-Tetrachloroethane | < 0.31 | 0.099 | 0.31 |
| Tetrachloroethene | 15 | 0.17 | 0.56 |
| Toluene | < 0.69 | 0.22 | 0.69 |
| 1,2,3-Trichlorobenzene | < 1 | 0.31 | 1.1 |
| 1,2,4-Trichlorobenzene | < 0.91 | 0.26 | 0.91 |
| 1,1,1-Trichloroethane | 0.97 | 0.2 | 0.63 |
| 1,1,2-Trichloroethane | < 0.17 | 0.055 | 0.17 |
| Trichloroethene | 4.7 | 0.055 | 0.18 |
| Trichlorofluoromethane | < 1.4 | 1.4 | 4.4 |
| 1,2,4-Trimethylbenzene | < 0.57 | 0.18 | 0.57 |
| 1,3,5-Trimethylbenzene | < 0.57 | 0.18 | 0.57 |
| Vinyl Chloride | 1.7 | 0.17 | 0.54 |
| m&p-Xylene | < 0.9 | 0.28 | 0.9 |
| o-Xylene | < 0.33 | 0.1 | 0.33 |

MDL = Method Detection Limit

PQL = Practical Quantitation Limit

NA = Not Applicable

Authorized Signature

**Analytical Laboratory**

1090 Kennedy Ave. Kimberly, WI 54136
414-735-8295

WI DNR Certified Lab #445027660

QC Summary**Method 8021 Volatile Organic Compounds**

Project #: CLW131246
Sample ID: MW200

Report Date: 02-Oct-96
Lab Code: 5014513A

| ANALYTE | INITIAL CALIBRATION | KNOWN STANDARD | MATRIX SPIKE | REPLICATE SPIKE | BLANK | PID SURROGATE | HALL SURROGATE |
|-----------------------------|---------------------|----------------|--------------|-----------------|-------|---------------|----------------|
| Benzene | P | P | P | P | P | P | P |
| Bromobenzene | P | P | P | P | P | P | P |
| Bromodichloromethane | P | P | P | P | P | P | P |
| n-Butylbenzene | P | P | P | P | P | P | P |
| sec-Butylbenzene | P | P | P | P | P | P | P |
| tert-Butylbenzene | P | P | P | P | P | P | P |
| Carbon Tetrachloride | P | P | P | P | P | P | P |
| Chlorobenzene | P | P | P | P | P | P | P |
| Chloroethane | P | F | P | P | P | P | P |
| Chloroform | P | P | F | P | F | P | P |
| Chloromethane | P | F | P | P | P | P | P |
| 2-Chlorotoluene | P | P | P | P | P | P | P |
| 4-Chlorotoluene | P | P | P | P | P | P | P |
| 1,2-Dibromo-3-Chloropropane | P | F | P | P | P | P | P |
| Dibromochloromethane | P | P | P | P | P | P | P |
| 1,2-Dichlorobenzene | P | P | P | P | P | P | P |
| 1,3-Dichlorobenzene | P | P | P | P | P | P | P |
| 1,4-Dichlorobenzene | P | P | P | P | P | P | P |
| Dichlorodifluoromethane | P | F | F | P | P | P | P |
| 1,1-Dichloroethane | P | P | P | P | P | P | P |
| 1,2-Dichloroethane | P | P | P | P | P | P | P |
| 1,1-Dichloroethene | P | F | P | P | P | P | P |
| cis-1,2-Dichloroethene | P | P | P | P | P | P | P |
| trans-1,2-Dichloroethene | P | P | P | P | P | P | P |
| 1,2-Dichloropropane | P | P | P | P | P | P | P |
| 1,3-Dichloropropane | P | P | F | P | P | P | P |
| 2,2-Dichloropropane | P | F | F | F | P | P | P |
| Di-Isopropyl Ether | P | P | P | P | P | P | P |
| Ethylbenzene | P | P | P | P | P | P | P |
| EDB (1,2-Dibromoethane) | P | P | P | P | P | P | P |
| Hexachlorobutadiene | P | P | P | P | P | P | P |
| Isopropylbenzene | P | P | P | P | P | P | P |
| p-Isopropyltoluene | P | P | P | P | P | P | P |
| Methylene Chloride | P | P | P | P | P | P | P |
| MTBE | P | P | P | P | P | P | P |
| Naphthalene | P | P | P | F | P | P | P |
| n-Propylbenzene | P | P | P | P | P | P | P |
| 1,1,2,2-Tetrachloroethane | P | P | P | F | P | P | P |
| Tetrachloroethene | P | P | P | P | P | P | P |
| Toluene | P | P | P | P | P | P | P |
| 1,2,3-Trichlorobenzene | P | P | P | P | P | P | P |
| 1,2,4-Trichlorobenzene | P | P | P | P | P | P | P |
| 1,1,1-Trichloroethane | P | P | P | P | P | P | P |
| 1,1,2-Trichloroethane | P | P | P | P | P | P | P |
| Trichloroethene | P | P | P | P | P | P | P |
| Trichlorofluoromethane | P | P | P | P | P | P | P |
| 1,2,4-Trimethylbenzene | P | P | P | P | P | P | P |
| 1,3,5-Trimethylbenzene | P | P | P | P | P | P | P |
| Vinyl Chloride | P | F | P | F | P | P | P |
| m & p-Xylene | P | P | P | P | P | P | P |
| o-Xylene | P | P | P | P | P | P | P |

P = Passed QC limits.

F = Failed QC limits.

NA = Not Applicable

VOC analysis detected unidentified peaks.

Authorized Signature

**Analytical Laboratory**

1090 Kennedy Ave. Kimberly, WI 54136
414-735-8295

WI DNR Certified Lab #445027660

BRIAN KRANZ
NORTHERN ENVIRONMENTAL
1214 W VENTURE COURT
MEQUON WI 53092

Project #: CLW131246
Project : Cedarburg
Sample ID: MW300
Lab Code: 5014513B
Sample Type: Water
Sample Date: 13-Sep-96

Report Date: 02-Oct-96

| Test | Result | MDL | PQL | Unit | pH | Date Ext/Digested | Date Analyzed: | Analyzed By: | QC Code |
|-----------------------------|--------|-----|-----|------|-----|-------------------|----------------|--------------|---------|
| MODIFIED DRO WDNR SEP 95 | < 100 | 30 | 96 | UG/L | 2.4 | 19-Sep-96 | 20-Sep-96 | C. Rotar | 1 |

MDL = Method Detection Limit

PQL = Practical Quantitation Limit

QC SUMMARY

CODE:

1 pH adjusted below two.

Authorized Signature

**Analytical Laboratory**

1090 Kennedy Ave. Kimberly, WI 54136
414-735-8295

WI DNR Certified Lab #445027660

Method 8021 Volatile Organic Compounds

BRIAN KRANZ
NORTHERN ENVIRONMENTAL
1214 W VENTURE COURT
MEQUON WI 53092

Report Date: 02-Oct-96
Analyzed By: G.Shah

Project #: CLW131246
Project : Cedarburg
Sample ID: MW300
Lab Code: 5014513B
Sample Type: Water
Sample Date: 13-Sep-96
Date Analyzed: 25-Sep-96

| ANALYTE | RESULT | MDL UG/L | PQL UG/L |
|-----------------------------|--------|-------------|-------------|
| Benzene | 0.34 | 0.082 | 0.26 |
| Bromobenzene | < 0.24 | 0.075 | 0.24 |
| Bromodichloromethane | < 0.11 | 0.035 | 0.11 |
| n-Butylbenzene | < 0.45 | 0.14 | 0.45 |
| sec-Butylbenzene | < 0.49 | 0.15 | 0.49 |
| tert-Butylbenzene | < 0.4 | 0.12 | 0.4 |
| Carbon Tetrachloride | < 0.5 | 0.16 | 0.5 |
| Chlorobenzene | < 0.27 | 0.086 | 0.27 |
| Chloroethane | 1.5 | 0.39 | 1.3 |
| Chloroform | < 0.22 | 0.07 | 0.22 |
| Chloromethane | < 1 | 0.88 | 3.1 |
| 2-Chlorotoluene | < 0.65 | 0.21 | 0.65 |
| 4-Chlorotoluene | < 0.19 | 0.06 | 0.19 |
| 1,2-Dibromo-3-Chloropropane | < 1 | 0.83 | 2.7 |
| Dibromochloromethane | < 0.09 | 0.028 | 0.09 |
| 1,2-Dichlorobenzene | 0.14 | 0.035 | 0.11 |
| 1,3-Dichlorobenzene | < 0.83 | 0.23 | 0.83 |
| 1,4-Dichlorobenzene | < 0.13 | 0.039 | 0.13 |
| Dichlorodifluoromethane | < 5.4 | 1.7 | 5.4 |
| 1,1-Dichloroethene | < 0.37 | 0.12 | 0.37 |
| 1,2-Dichloroethane | < 0.86 | 0.27 | 0.86 |
| 1,1-Dichloroethane | 0.63 | 0.084 | 0.27 |
| cis 1,2-Dichloroethene | 0.59 | 0.092 | 0.29 |
| trans-1,2-dichloroethene | < 0.23 | 0.072 | 0.23 |
| 1,2-Dichloropropane | < 0.15 | 0.046 | 0.15 |
| 1,3-DCP, Tetrachloroethene | < 0.56 | 0.17 | 0.56 |

Fluorobenzene Surrogate 108 % Rec.
1,4-Dichlorobutane Surrogate 87 % Rec.
Sample pH 1.6

| ANALYTE | RESULT | MDL UG/L | PQL UG/L |
|---------------------------|--------|-------------|-------------|
| 2,2-Dichloropropane | < 1 | 0.63 | 2.2 |
| Di-isopropyl Ether | < 0.38 | 0.12 | 0.38 |
| Ethylbenzene | < 0.32 | 0.1 | 0.32 |
| EDB (1,2-Dibromoethane) | < 0.08 | 0.025 | 0.08 |
| Hexachlorobutadiene | < 0.35 | 0.11 | 0.35 |
| Isopropylbenzene | < 0.36 | 0.11 | 0.36 |
| p-Isopropyltoluene | < 0.46 | 0.15 | 0.46 |
| Methylene Chloride | < 4 | 0.29 | 0.91 |
| MTBE | < 0.22 | 0.069 | 0.22 |
| Naphthalene | 0.56 | 0.13 | 0.41 |
| n-Propylbenzene | < 0.41 | 0.13 | 0.41 |
| 1,1,2,2-Tetrachloroethane | < 0.31 | 0.099 | 0.31 |
| Tetrachloroethene | 3.2 | 0.17 | 0.56 |
| Toluene | < 0.69 | 0.22 | 0.69 |
| 1,2,3-Trichlorobenzene | < 1 | 0.31 | 1.1 |
| 1,2,4-Trichlorobenzene | < 0.91 | 0.26 | 0.91 |
| 1,1,1-Trichloroethane | < 0.63 | 0.2 | 0.63 |
| 1,1,2-Trichloroethane | < 0.17 | 0.055 | 0.17 |
| Trichloroethene | 0.49 | 0.055 | 0.18 |
| Trichlorofluoromethane | < 1.4 | 1.4 | 4.4 |
| 1,2,4-Trimethylbenzene | < 0.57 | 0.18 | 0.57 |
| 1,3,5-Trimethylbenzene | < 0.57 | 0.18 | 0.57 |
| Vinyl Chloride | < 0.54 | 0.17 | 0.54 |
| m&p-Xylene | < 0.9 | 0.28 | 0.9 |
| o-Xylene | < 0.33 | 0.1 | 0.33 |

MDL = Method Detection Limit

PQL = Practical Quantitation Limit

NA = Not Applicable

Authorized Signature

**Analytical Laboratory**

1090 Kennedy Ave. Kimberly, WI 54136
414-735-8295

WI DNR Certified Lab #445027660

QC Summary**Method 8021 Volatile Organic Compounds**Project #: CLW131246
Sample ID:

MW300

Report Date: 02-Oct-96
Lab Code: 5014513B

| ANALYTE | INITIAL CALIBRATION | KNOWN STANDARD | MATRIX SPIKE | REPLICATE SPIKE | BLANK | PID SURROGATE | HALL SURROGATE |
|-----------------------------|---------------------|----------------|--------------|-----------------|-------|---------------|----------------|
| Benzene | P | P | P | P | P | P | P |
| Bromobenzene | P | P | P | P | P | P | P |
| Bromodichloromethane | P | P | P | P | P | P | P |
| n-Butylbenzene | P | P | P | P | P | P | P |
| sec-Butylbenzene | P | P | P | P | P | P | P |
| tert-Butylbenzene | P | P | P | P | P | P | P |
| Carbon Tetrachloride | P | P | P | P | P | P | P |
| Chlorobenzene | P | P | P | P | P | P | P |
| Chloroethane | P | F | P | P | P | P | P |
| Chloroform | P | P | P | P | P | P | P |
| Chloromethane | P | F | P | P | F | P | P |
| 2-Chlorotoluene | P | P | P | P | P | P | P |
| 4-Chlorotoluene | P | P | P | P | P | P | P |
| 1,2-Dibromo-3-Chloropropane | P | F | P | P | P | P | P |
| Dibromochloromethane | P | P | P | P | P | P | P |
| 1,2-Dichlorobenzene | P | P | P | P | P | P | P |
| 1,3-Dichlorobenzene | P | P | P | P | P | P | P |
| 1,4-Dichlorobenzene | P | F | P | P | P | P | P |
| Dichlorodifluoromethane | P | P | F | P | P | P | P |
| 1,1-Dichloroethane | P | P | P | P | P | P | P |
| 1,2-Dichloroethane | P | P | P | P | P | P | P |
| 1,1-Dichloroethene | P | F | P | P | P | P | P |
| trans-1,2-Dichloroethene | P | P | P | P | P | P | P |
| 1,2-Dichloropropane | P | P | P | P | P | P | P |
| 1,3-Dichloropropane | P | F | P | F | P | P | P |
| 2,2-Dichloropropane | P | P | F | P | P | P | P |
| Di-Isopropyl Ether | P | P | P | P | P | P | P |
| Ethylbenzene | P | P | P | P | P | P | P |
| EDB (1,2-Dibromoethane) | P | P | P | P | P | P | P |
| Hexachlorobutadiene | P | P | P | P | P | P | P |
| Isopropylbenzene | P | P | P | P | P | P | P |
| p-Isopropyltoluene | P | P | P | P | P | P | P |
| Methylene Chloride | P | P | P | P | P | P | P |
| MTBE | P | P | P | P | P | P | P |
| Naphthalene | P | P | P | F | P | P | P |
| n-Propylbenzene | P | P | P | P | P | P | P |
| 1,1,2,2-Tetrachloroethane | P | P | P | F | P | P | P |
| Tetrachloroethene | P | P | P | P | P | P | P |
| Toluene | P | P | P | P | P | P | P |
| 1,2,3-Trichlorobenzene | P | P | P | P | P | P | P |
| 1,2,4-Trichlorobenzene | P | P | P | P | P | P | P |
| 1,1,1-Trichloroethane | P | P | P | P | P | P | P |
| 1,1,2-Trichloroethane | P | P | P | P | P | P | P |
| Trichloroethene | P | P | P | P | P | P | P |
| Trichlorofluoromethane | P | P | P | P | P | P | P |
| 1,2,4-Trimethylbenzene | P | P | P | P | P | P | P |
| 1,3,5-Trimethylbenzene | P | P | P | P | P | P | P |
| Vinyl Chloride | P | F | P | F | P | P | P |
| m & p-Xylene | P | P | P | P | P | P | P |
| o-Xylene | P | P | P | P | P | P | P |

P = Passed QC limits.

F = Failed QC limits.

NA = Not Applicable

VOC analysis detected unidentified peaks.

Authorized Signature

**Analytical Laboratory**

1090 Kennedy Ave. Kimberly, WI 54136
414-735-8295

WI DNR Certified Lab #445027660

BRIAN KRANZ
NORTHERN ENVIRONMENTAL
1214 W VENTURE COURT
MEQUON WI 53092

Project #: CLW131246
Project : Cedarburg
Sample ID: MW400
Lab Code: 5014513C
Sample Type: Water
Sample Date: 13-Sep-96

Report Date: 02-Oct-96

| Test | Result | MDL | PQL | Unit | pH | Date Ext/Digested | Date Analyzed: | Analyzed By: | QC Code |
|-----------------------------|--------|-----|-----|------|-----|-------------------|----------------|--------------|---------|
| MODIFIED DRO WDNR SEP 95 | < 100 | 30 | 96 | UG/L | 2.5 | 19-Sep-96 | 20-Sep-96 | C. Rotar | 1 |

MDL = Method Detection Limit

PQL = Practical Quantitation Limit

QC SUMMARY

CODE:

1 pH adjusted below two.

Authorized Signature

**Analytical Laboratory**

1090 Kennedy Ave. Kimberly, WI 54136
414-735-8295

WI DNR Certified Lab #445027660

Method 8021 Volatile Organic Compounds

BRIAN KRANZ
NORTHERN ENVIRONMENTAL
1214 W VENTURE COURT
MEQUON WI 53092

Report Date: 24-Sep-96
Analyzed By: R. Everson

Project #: CLW131246
Project : Cedarburg
Sample ID: MW400
Lab Code: 5014513C
Sample Type: Water
Sample Date: 13-Sep-96
Date Analyzed: 19-Sep-96

| ANALYTE | RESULT | MDL UG/L | PQL UG/L |
|-----------------------------|--------|-------------|-------------|
| Benzene | < 0.26 | 0.082 | 0.26 |
| Bromobenzene | < 0.24 | 0.075 | 0.24 |
| Bromodichloromethane | < 0.11 | 0.035 | 0.11 |
| n-Butylbenzene | < 0.45 | 0.14 | 0.45 |
| sec-Butylbenzene | < 0.49 | 0.15 | 0.49 |
| tert-Butylbenzene | < 0.4 | 0.12 | 0.4 |
| Carbon Tetrachloride | < 0.5 | 0.16 | 0.5 |
| Chlorobenzene | < 0.27 | 0.086 | 0.27 |
| Chloroethane | < 0.5 | 0.39 | 1.3 |
| Chloroform | < 0.22 | 0.07 | 0.22 |
| Chloromethane | < 1 | 0.88 | 3.1 |
| 2-Chlorotoluene | < 0.65 | 0.21 | 0.65 |
| 4-Chlorotoluene | < 0.19 | 0.06 | 0.19 |
| 1,2-Dibromo-3-Chloropropane | < 1 | 0.83 | 2.7 |
| Dibromochloromethane | < 0.09 | 0.028 | 0.09 |
| 1,2-Dichlorobenzene | < 0.11 | 0.035 | 0.11 |
| 1,3-Dichlorobenzene | < 0.83 | 0.23 | 0.83 |
| 1,4-Dichlorobenzene | < 0.13 | 0.039 | 0.13 |
| Dichlorodifluoromethane | < 5.4 | 1.7 | 5.4 |
| 1,1-Dichloroethene | < 0.37 | 0.12 | 0.37 |
| 1,2-Dichloroethane | < 0.86 | 0.27 | 0.86 |
| 1,1-Dichloroethane | < 0.27 | 0.084 | 0.27 |
| cis 1,2-Dichloroethene | < 0.29 | 0.092 | 0.29 |
| trans-1,2-dichloroethene | < 0.23 | 0.072 | 0.23 |
| 1,2-Dichloropropane | < 0.15 | 0.046 | 0.15 |
| 1,3-DCP, Tetrachloroethene | < 0.56 | 0.17 | 0.56 |

Fluorobenzene Surrogate 108 % Rec.
1,4-Dichlorobutane Surrogate 89 % Rec.
Sample pH 1.6

| ANALYTE | RESULT | MDL UG/L | PQL UG/L |
|---------------------------|--------|-------------|-------------|
| 2,2-Dichloropropane | < 1 | 0.63 | 2.2 |
| Di-isopropyl Ether | < 0.38 | 0.12 | 0.38 |
| Ethylbenzene | < 0.32 | 0.1 | 0.32 |
| EDB (1,2-Dibromoethane) | < 0.08 | 0.025 | 0.08 |
| Hexachlorobutadiene | < 0.35 | 0.11 | 0.35 |
| Isopropylbenzene | < 0.36 | 0.11 | 0.36 |
| p-Isopropyltoluene | < 0.46 | 0.15 | 0.46 |
| Methylene Chloride | < 4 | 0.29 | 0.91 |
| MTBE | < 0.22 | 0.069 | 0.22 |
| Naphthalene | < 0.41 | 0.13 | 0.41 |
| n-Propylbenzene | < 0.41 | 0.13 | 0.41 |
| 1,1,2,2-Tetrachloroethane | < 0.31 | 0.099 | 0.31 |
| Tetrachloroethene | 0.68 | 0.17 | 0.56 |
| Toluene | < 0.69 | 0.22 | 0.69 |
| 1,2,3-Trichlorobenzene | < 1 | 0.31 | 1.1 |
| 1,2,4-Trichlorobenzene | < 0.91 | 0.26 | 0.91 |
| 1,1,1-Trichloroethane | < 0.63 | 0.2 | 0.63 |
| 1,1,2-Trichloroethane | < 0.17 | 0.055 | 0.17 |
| Trichloroethene | < 0.18 | 0.055 | 0.18 |
| Trichlorofluoromethane | < 1.4 | 1.4 | 4.4 |
| 1,2,4-Trimethylbenzene | < 0.57 | 0.18 | 0.57 |
| 1,3,5-Trimethylbenzene | < 0.57 | 0.18 | 0.57 |
| Vinyl Chloride | < 0.54 | 0.17 | 0.54 |
| m&p-Xylene | < 0.9 | 0.28 | 0.9 |
| o-Xylene | < 0.33 | 0.1 | 0.33 |

MDL = Method Detection Limit

PQL = Practical Quantitation Limit

NA = Not Applicable

Authorized Signature

**Analytical Laboratory**

1090 Kennedy Ave. Kimberly, WI 54136
414-735-8295

WI DNR Certified Lab #445027660**QC Summary****Method 8021 Volatile Organic Compounds**

Project #: CLW131246 Report Date: 24-Sep-96
Sample ID: MW400 Lab Code: 5014513C

| ANALYTE | INITIAL CALIBRATION | KNOWN STANDARD | MATRIX SPIKE | REPLICATE SPIKE | BLANK | PID SURROGATE | HALL SURROGATE |
|-----------------------------|---------------------|----------------|--------------|-----------------|-------|---------------|----------------|
| Benzene | P | P | P | P | P | P | P |
| Bromobenzene | P | P | P | P | P | P | P |
| Bromodichloromethane | P | P | P | F | P | P | P |
| n-Butylbenzene | P | P | P | P | P | P | P |
| sec-Butylbenzene | P | P | P | P | P | P | P |
| tert-Butylbenzene | P | P | P | P | P | P | P |
| Carbon Tetrachloride | P | P | P | P | P | P | P |
| Chlorobenzene | P | P | P | P | P | P | P |
| Chloroethane | P | P | P | P | P | P | P |
| Chloroform | P | P | P | P | P | P | P |
| Chloromethane | P | F | P | P | P | P | P |
| 2-Chlorotoluene | P | P | P | F | P | P | P |
| 4-Chlorotoluene | P | P | P | F | P | P | P |
| 1,2-Dibromo-3-Chloropropane | P | F | P | F | P | P | P |
| Dibromochloromethane | P | P | P | F | P | P | P |
| 1,2-Dichlorobenzene | P | P | P | P | P | P | P |
| 1,3-Dichlorobenzene | P | P | P | P | P | P | P |
| 1,4-Dichlorobenzene | P | P | P | F | P | P | P |
| Dichlorodifluoromethane | P | F | P | P | P | P | P |
| 1,1-Dichloroethane | P | P | P | P | P | P | P |
| 1,2-Dichloroethane | P | P | P | F | P | P | P |
| 1,1-Dichloroethene | P | F | P | P | P | P | P |
| cis-1,2-Dichloroethene | P | F | P | P | P | P | P |
| trans-1,2-Dichloroethene | P | F | P | P | P | P | P |
| 1,2-Dichloropropane | P | P | P | F | P | P | P |
| 1,3-Dichloropropane | P | P | P | F | P | P | P |
| 2,2-Dichloropropane | P | F | P | F | P | P | P |
| Di-Isopropyl Ether | P | P | P | P | P | P | P |
| Ethylbenzene | P | P | P | P | P | P | P |
| EDB (1,2-Dibromoethane) | P | P | P | F | P | P | P |
| Hexachlorobutadiene | P | P | P | F | P | P | P |
| Isopropylbenzene | P | P | P | P | P | P | P |
| p-Isopropyltoluene | P | P | P | P | P | P | P |
| Methylene Chloride | P | F | P | P | P | P | P |
| MTBE | P | P | P | P | P | P | P |
| Naphthalene | P | P | P | P | P | P | P |
| n-Propylbenzene | P | P | P | P | P | P | P |
| 1,1,2,2-Tetrachloroethane | P | P | P | F | P | P | P |
| Tetrachloroethene | P | P | P | P | P | P | P |
| Toluene | P | P | P | P | P | P | P |
| 1,2,3-Trichlorobenzene | P | P | P | F | P | P | P |
| 1,2,4-Trichlorobenzene | P | P | P | F | P | P | P |
| 1,1,1-Trichloroethane | P | P | P | P | P | P | P |
| 1,1,2-Trichloroethane | P | P | P | F | P | P | P |
| Trichloroethene | P | P | P | P | P | P | P |
| Trichlorofluoromethane | P | F | F | P | P | P | P |
| 1,2,4-Trimethylbenzene | P | P | P | P | P | P | P |
| 1,3,5-Trimethylbenzene | P | F | P | P | P | P | P |
| Vinyl Chloride | P | P | P | P | P | P | P |
| m & p-Xylene | P | F | P | P | P | P | P |
| o-Xylene | P | P | P | P | P | P | P |

P = Passed QC limits.

F = Failed QC limits.

NA = Not Applicable

VOC analysis detected unidentified peaks.

Authorized Signature

**Analytical Laboratory**

1090 Kennedy Ave. Kimberly, WI 54136
414-735-8295

WI DNR Certified Lab #445027660

BRIAN KRANZ
NORTHERN ENVIRONMENTAL
1214 W VENTURE COURT
MEQUON WI 53092

Project #: CLW131246
Project : Cedarburg
Sample ID: MW500
Lab Code: 5014513D
Sample Type: Water
Sample Date: 13-Sep-96

Report Date: 02-Oct-96

| Test | Result | MDL | PQL | Unit | pH | Date Ext/Digested | Date Analyzed: | Analyzed By: | QC Code |
|-----------------------------|--------|-----|-----|------|-----|-------------------|----------------|--------------|---------|
| MODIFIED DRO WDNR SEP 95 | < 100 | 30 | 96 | UG/L | 2.5 | 19-Sep-96 | 20-Sep-96 | C. Rotar | 1 |

MDL = Method Detection Limit

PQL = Practical Quantitation Limit

QC SUMMARY

CODE:

1 pH adjusted below two.

Authorized Signature


Analytical Laboratory

 1090 Kennedy Ave. Kimberly, WI 54136
 414-735-8295

WI DNR Certified Lab #445027660

Method 8021 Volatile Organic Compounds

BRIAN KRANZ
 NORTHERN ENVIRONMENTAL
 1214 W VENTURE COURT
 MEQUON WI 53092

Report Date: 24-Sep-96
 Analyzed By: R. Everson

Project #: CLW131246
 Project : Cedarburg
 Sample ID: MW500
 Lab Code: 5014513D
 Sample Type: Water
 Sample Date: 13-Sep-96
 Date Analyzed: 19-Sep-96

| ANALYTE | RESULT | MDL UG/L | PQL UG/L |
|-----------------------------|--------|-------------|-------------|
| Benzene | < 0.26 | 0.082 | 0.26 |
| Bromobenzene | < 0.24 | 0.075 | 0.24 |
| Bromodichloromethane | < 0.11 | 0.035 | 0.11 |
| n-Butylbenzene | < 0.45 | 0.14 | 0.45 |
| sec-Butylbenzene | < 0.49 | 0.15 | 0.49 |
| tert-Butylbenzene | < 0.4 | 0.12 | 0.4 |
| Carbon Tetrachloride | < 0.5 | 0.16 | 0.5 |
| Chlorobenzene | < 0.27 | 0.086 | 0.27 |
| Chloroethane | < 0.5 | 0.39 | 1.3 |
| Chloroform | < 0.22 | 0.07 | 0.22 |
| Chloromethane | < 1 | 0.88 | 3.1 |
| 2-Chlorotoluene | < 0.65 | 0.21 | 0.65 |
| 4-Chlorotoluene | < 0.19 | 0.06 | 0.19 |
| 1,2-Dibromo-3-Chloropropane | < 1 | 0.83 | 2.7 |
| Dibromochloromethane | < 0.09 | 0.028 | 0.09 |
| 1,2-Dichlorobenzene | < 0.11 | 0.035 | 0.11 |
| 1,3-Dichlorobenzene | < 0.83 | 0.23 | 0.83 |
| 1,4-Dichlorobenzene | < 0.13 | 0.039 | 0.13 |
| Dichlorodifluoromethane | < 5.4 | 1.7 | 5.4 |
| 1,1-Dichloroethene | < 0.37 | 0.12 | 0.37 |
| 1,2-Dichloroethane | < 0.86 | 0.27 | 0.86 |
| 1,1-Dichloroethane | < 0.27 | 0.084 | 0.27 |
| cis 1,2-Dichloroethene | < 0.29 | 0.092 | 0.29 |
| trans-1,2-dichloroethene | < 0.23 | 0.072 | 0.23 |
| 1,2-Dichloropropane | < 0.15 | 0.046 | 0.15 |
| 1,3-DCP, Tetrachloroethene | < 0.56 | 0.17 | 0.56 |

| ANALYTE | RESULT | MDL UG/L | PQL UG/L |
|---------------------------|--------|-------------|-------------|
| 2,2-Dichloropropane | < 1 | 0.63 | 2.2 |
| Di-isopropyl Ether | < 0.38 | 0.12 | 0.38 |
| Ethylbenzene | < 0.32 | 0.1 | 0.32 |
| EDB (1,2-Dibromoethane) | < 0.08 | 0.025 | 0.08 |
| Hexachlorobutadiene | < 0.35 | 0.11 | 0.35 |
| Isopropylbenzene | < 0.36 | 0.11 | 0.36 |
| p-Isopropyltoluene | < 0.46 | 0.15 | 0.46 |
| Methylene Chloride | < 4 | 0.29 | 0.91 |
| MTBE | < 0.22 | 0.069 | 0.22 |
| Naphthalene | < 0.41 | 0.13 | 0.41 |
| n-Propylbenzene | < 0.41 | 0.13 | 0.41 |
| 1,1,2,2-Tetrachloroethane | < 0.31 | 0.099 | 0.31 |
| Tetrachloroethene | < 0.56 | 0.17 | 0.56 |
| Toluene | < 0.69 | 0.22 | 0.69 |
| 1,2,3-Trichlorobenzene | < 1 | 0.31 | 1.1 |
| 1,2,4-Trichlorobenzene | < 0.91 | 0.26 | 0.91 |
| 1,1,1-Trichloroethane | < 0.63 | 0.2 | 0.63 |
| 1,1,2-Trichloroethane | < 0.17 | 0.055 | 0.17 |
| Trichloroethene | < 0.18 | 0.055 | 0.18 |
| Trichlorofluoromethane | < 1.4 | 1.4 | 4.4 |
| 1,2,4-Trimethylbenzene | < 0.57 | 0.18 | 0.57 |
| 1,3,5-Trimethylbenzene | < 0.57 | 0.18 | 0.57 |
| Vinyl Chloride | < 0.54 | 0.17 | 0.54 |
| m&p-Xylene | < 0.9 | 0.28 | 0.9 |
| o-Xylene | < 0.33 | 0.1 | 0.33 |

Fluorobenzene Surrogate 108 % Rec.
 1,4-Dichlorobutane Surrogate 86 % Rec.
 Sample pH 1.6

MDL = Method Detection Limit

PQL = Practical Quantitation Limit

NA = Not Applicable

Authorized Signature

**Analytical Laboratory**

1090 Kennedy Ave. Kimberly, WI 54136
414-735-8295

WI DNR Certified Lab #445027660

QC Summary**Method 8021 Volatile Organic Compounds**

Project #: CLW131246 Report Date: 24-Sep-96
Sample ID: MW500 Lab Code: 5014513D

| ANALYTE | INITIAL CALIBRATION | KNOWN STANDARD | MATRIX SPIKE | REPLICATE SPIKE | BLANK | PID SURROGATE | HALL SURROGATE |
|-----------------------------|---------------------|----------------|--------------|-----------------|-------|---------------|----------------|
| Benzene | P | P | P | P | P | P | P |
| Bromobenzene | P | P | P | P | P | P | P |
| Bromodichloromethane | P | P | P | F | P | P | P |
| n-Butylbenzene | P | P | P | P | P | P | P |
| sec-Butylbenzene | P | P | P | P | P | P | P |
| tert-Butylbenzene | P | P | P | P | P | P | P |
| Carbon Tetrachloride | P | P | P | P | P | P | P |
| Chlorobenzene | P | P | P | P | P | P | P |
| Chloroethane | P | P | P | P | P | P | P |
| Chloroform | P | P | P | P | P | P | P |
| Chloromethane | P | F | P | P | P | P | P |
| 2-Chlorotoluene | P | P | P | F | P | P | P |
| 4-Chlorotoluene | P | P | P | F | P | P | P |
| 1,2-Dibromo-3-Chloropropane | P | F | P | F | P | P | P |
| Dibromochloromethane | P | P | P | F | P | P | P |
| 1,2-Dichlorobenzene | P | P | P | P | P | P | P |
| 1,3-Dichlorobenzene | P | P | P | P | P | P | P |
| 1,4-Dichlorobenzene | P | P | P | F | P | P | P |
| Dichlorodifluoromethane | P | F | P | P | P | P | P |
| 1,1-Dichloroethane | P | P | P | P | P | P | P |
| 1,2-Dichloroethane | P | P | P | F | P | P | P |
| 1,1-Dichloroethene | P | F | P | P | P | P | P |
| cis-1,2-Dichloroethene | P | P | P | P | P | P | P |
| trans-1,2-Dichloroethene | P | F | P | P | P | P | P |
| 1,2-Dichloropropene | P | P | P | F | P | P | P |
| 1,3-Dichloropropene | P | P | P | F | P | P | P |
| 2,2-Dichloropropene | P | F | P | F | P | P | P |
| Di-Isopropyl Ether | P | P | P | P | P | P | P |
| Ethylbenzene | P | P | P | P | P | P | P |
| EDB (1,2-Dibromoethane) | P | P | P | P | F | P | P |
| Hexachlorobutadiene | P | P | P | P | F | P | P |
| Isopropylbenzene | P | P | P | P | P | P | P |
| p-Isopropyltoluene | P | P | P | P | P | P | P |
| Methylene Chloride | P | F | P | P | P | P | P |
| MTBE | P | P | P | P | P | P | P |
| Naphthalene | P | P | P | P | P | P | P |
| n-Propylbenzene | P | P | P | P | P | P | P |
| 1,1,2-Tetrachloroethane | P | P | P | P | F | P | P |
| Tetrachloroethene | P | P | P | P | P | P | P |
| Toluene | P | P | P | P | P | P | P |
| 1,2,3-Trichlorobenzene | P | P | P | P | F | P | P |
| 1,2,4-Trichlorobenzene | P | P | P | P | F | P | P |
| 1,1,1-Trichloroethane | P | P | P | P | F | P | P |
| 1,1,2-Trichloroethane | P | P | P | P | F | P | P |
| Trichloroethene | P | P | P | P | P | P | P |
| Trichlorofluoromethane | P | F | P | P | P | P | P |
| 1,2,4-Trimethylbenzene | P | P | P | P | P | P | P |
| 1,3,5-Trimethylbenzene | P | F | P | P | P | P | P |
| Vinyl Chloride | P | F | P | P | P | P | P |
| m & p-Xylene | P | P | P | P | P | P | P |
| o-Xylene | P | P | P | P | P | P | P |

P = Passed QC limits.

F = Failed QC limits.

NA = Not Applicable

VOC analysis detected unidentified peaks.

Authorized Signature


Analytical Laboratory

1090 Kennedy Ave. Kimberly, WI 54136
414-735-8295

WI DNR Certified Lab #445027660

Method 8021 Volatile Organic Compounds

BRIAN KRANZ
NORTHERN ENVIRONMENTAL
1214 W VENTURE COURT
MEQUON WI 53092

Report Date: 03-Oct-96
Analyzed By: G.Shah

Project #: CLW131246
Project : Cedarburg
Sample ID: Dup-1246
Lab Code: 5014513E
Sample Type: Water
Sample Date: 13-Sep-96
Date Analyzed: 26-Sep-96

| ANALYTE | RESULT | MDL UG/L | PQL UG/L |
|-----------------------------|--------|-------------|-------------|
| Benzene | 0.52 | 0.082 | 0.26 |
| Bromobenzene | < 0.24 | 0.075 | 0.24 |
| Bromodichloromethane | < 0.11 | 0.035 | 0.11 |
| n-Butylbenzene | < 0.45 | 0.14 | 0.45 |
| sec-Butylbenzene | 0.7 | 0.15 | 0.49 |
| tert-Butylbenzene | < 0.4 | 0.12 | 0.4 |
| Carbon Tetrachloride | < 0.5 | 0.16 | 0.5 |
| Chlorobenzene | < 0.27 | 0.086 | 0.27 |
| Chloroethane | 2.6 | 0.39 | 1.3 |
| Chloroform | < 0.22 | 0.07 | 0.22 |
| Chloromethane | < 1 | 0.88 | 3.1 |
| 2-Chlorotoluene | < 0.65 | 0.21 | 0.65 |
| 4-Chlorotoluene | < 0.19 | 0.06 | 0.19 |
| 1,2-Dibromo-3-Chloropropane | < 1 | 0.83 | 2.7 |
| Dibromochloromethane | < 0.09 | 0.028 | 0.09 |
| 1,2-Dichlorobenzene | 1 | 0.035 | 0.11 |
| 1,3-Dichlorobenzene | < 0.83 | 0.23 | 0.83 |
| 1,4-Dichlorobenzene | 0.14 | 0.039 | 0.13 |
| Dichlorodifluoromethane | < 5.4 | 1.7 | 5.4 |
| 1,1-Dichloroethene | < 0.37 | 0.12 | 0.37 |
| 1,2-Dichloroethane | < 0.86 | 0.27 | 0.86 |
| 1,1-Dichloroethane | 4.9 | 0.084 | 0.27 |
| cis 1,2-Dichloroethene | 4.1 | 0.092 | 0.29 |
| trans-1,2-dichloroethene | 0.3 | 0.072 | 0.23 |
| 1,2-Dichloropropane | < 0.15 | 0.046 | 0.15 |
| 1,3-DCP, Tetrachloroethene | < 0.56 | 0.17 | 0.56 |

| ANALYTE | RESULT | MDL UG/L | PQL UG/L |
|---------------------------|--------|-------------|-------------|
| 2,2-Dichloropropane | < 1 | 0.63 | 2.2 |
| Di-isopropyl Ether | < 0.38 | 0.12 | 0.38 |
| Ethylbenzene | < 0.32 | 0.1 | 0.32 |
| EDB (1,2-Dibromoethane) | 0.12 | 0.025 | 0.08 |
| Hexachlorobutadiene | < 0.35 | 0.11 | 0.35 |
| Isopropylbenzene | < 0.36 | 0.11 | 0.36 |
| p-Isopropyltoluene | < 0.46 | 0.15 | 0.46 |
| Methylene Chloride | < 4 | 0.29 | 0.91 |
| MTBE | < 0.22 | 0.069 | 0.22 |
| Naphthalene | < 0.41 | 0.13 | 0.41 |
| n-Propylbenzene | < 0.41 | 0.13 | 0.41 |
| 1,1,2,2-Tetrachloroethane | < 0.31 | 0.099 | 0.31 |
| Tetrachloroethene | 15 | 0.17 | 0.56 |
| Toluene | < 0.69 | 0.22 | 0.69 |
| 1,2,3-Trichlorobenzene | < 1 | 0.31 | 1.1 |
| 1,2,4-Trichlorobenzene | < 0.91 | 0.26 | 0.91 |
| 1,1,1-Trichloroethane | 1 | 0.2 | 0.63 |
| 1,1,2-Trichloroethane | < 0.17 | 0.055 | 0.17 |
| Trichloroethene | 4.8 | 0.055 | 0.18 |
| Trichlorofluoromethane | < 1.4 | 1.4 | 4.4 |
| 1,2,4-Trimethylbenzene | < 0.57 | 0.18 | 0.57 |
| 1,3,5-Trimethylbenzene | < 0.57 | 0.18 | 0.57 |
| Vinyl Chloride | 1.8 | 0.17 | 0.54 |
| m&p-Xylene | < 0.9 | 0.28 | 0.9 |
| o-Xylene | < 0.33 | 0.1 | 0.33 |

MDL = Method Detection Limit

PQL = Practical Quantitation Limit

NA = Not Applicable

Fluorobenzene Surrogate 106 % Rec.
1,4-Dichlorobutane Surrogate 87 % Rec.
Sample pH 1.6

Authorized Signature

**Analytical Laboratory**

1090 Kennedy Ave. Kimberly, WI 54136
414-735-8295

WI DNR Certified Lab #445027660

QC Summary**Method 8021 Volatile Organic Compounds**

Project #: CLW131246
Sample ID: Dup-1246

Report Date: 02-Oct-96
Lab Code: 5014513E

| ANALYTE | INITIAL CALIBRATION | KNOWN STANDARD | MATRIX SPIKE | REPLICATE SPIKE | BLANK | PID SURROGATE | HALL SURROGATE |
|-----------------------------|---------------------|----------------|--------------|-----------------|-------|---------------|----------------|
| Benzene | P | P | P | P | P | P | P |
| Bromobenzene | P | P | P | P | P | P | P |
| Bromodichloromethane | P | P | P | P | P | P | P |
| n-Butylbenzene | P | P | P | P | P | P | P |
| sec-Butylbenzene | P | P | P | P | P | P | P |
| tert-Butylbenzene | P | P | P | P | P | P | P |
| Carbon Tetrachloride | P | P | P | P | P | P | P |
| Chlorobenzene | P | P | P | P | P | P | P |
| Chloroethane | P | F | P | P | P | P | P |
| Chloroform | P | P | P | P | P | P | P |
| Chlormethane | P | F | F | P | F | P | P |
| 2-Chlorotoluene | P | P | P | P | P | P | P |
| 4-Chlorotoluene | P | P | P | P | P | P | P |
| 1,2-Dibromo-3-Chloropropane | P | F | P | P | P | P | P |
| Dibromochloromethane | P | P | P | P | P | P | P |
| 1,2-Dichlorobenzene | P | P | P | P | P | P | P |
| 1,3-Dichlorobenzene | P | P | P | P | P | P | P |
| 1,4-Dichlorobenzene | P | P | P | P | P | P | P |
| Dichlorodifluoromethane | P | F | F | P | P | P | P |
| 1,1-Dichloroethane | P | P | P | P | P | P | P |
| 1,2-Dichloroethane | P | P | P | P | P | P | P |
| 1,1-Dichloroethene | P | F | P | P | P | P | P |
| cis-1,2-Dichloroethene | P | P | P | P | P | P | P |
| trans-1,2-Dichloroethene | P | P | P | P | P | P | P |
| 1,2-Dichloropropane | P | P | P | P | P | P | P |
| 1,3-Dichloropropane | P | F | F | P | P | P | P |
| 2,2-Dichloropropane | P | F | F | F | P | P | P |
| Di-Isopropyl Ether | P | P | P | P | P | P | P |
| Ethylbenzene | P | P | P | P | P | P | P |
| EDB (1,2-Dibromoethane) | P | P | P | P | P | P | P |
| Hexachlorobutadiene | P | P | P | P | P | P | P |
| Isopropylbenzene | P | P | P | P | P | P | P |
| p-Isopropyltoluene | P | P | P | P | P | P | P |
| Methylene Chloride | P | P | P | P | P | P | P |
| MTBE | P | P | P | P | P | P | P |
| Naphthalene | P | P | P | F | P | P | P |
| n-Propylbenzene | P | P | P | P | P | P | P |
| 1,1,2,2-Tetrachloroethane | P | P | P | F | P | P | P |
| Tetrachloroethene | P | P | P | P | P | P | P |
| Toluene | P | P | P | P | P | P | P |
| 1,2,3-Trichlorobenzene | P | P | P | P | P | P | P |
| 1,2,4-Trichlorobenzene | P | P | P | P | P | P | P |
| 1,1,1-Trichloroethane | P | P | P | P | P | P | P |
| 1,1,2-Trichloroethane | P | P | P | P | P | P | P |
| Trichloroethene | P | P | P | P | P | P | P |
| Trichlorofluoromethane | P | P | P | P | P | P | P |
| 1,2,4-Trimethylbenzene | P | P | P | P | P | P | P |
| 1,3,5-Trimethylbenzene | P | F | P | P | P | P | P |
| Vinyl Chloride | P | P | P | P | P | P | P |
| m & p-Xylene | P | F | P | F | P | P | P |
| o-Xylene | P | P | P | P | P | P | P |

P = Passed QC limits.

F = Failed QC limits.

NA = Not Applicable

VOC analysis detected unidentified peaks.

Authorized Signature

**Analytical Laboratory**

1090 Kennedy Ave. Kimberly, WI 54136
414-735-8295

WI DNR Certified Lab #445027660

Method 8021 Volatile Organic Compounds

BRIAN KRANZ
NORTHERN ENVIRONMENTAL
1214 W VENTURE COURT
MEQUON WI 53092

Report Date: 24-Sep-96
Analyzed By: R. Everson

Project #: CLW131246
Project : Cedarburg
Sample ID: Trip
Lab Code: 5014513F
Sample Type: Water
Sample Date: 13-Sep-96
Date Analyzed: 18-Sep-96

| ANALYTE | RESULT | MDL UG/L | PQL UG/L |
|-----------------------------|--------|-------------|-------------|
| Benzene | < 0.26 | 0.082 | 0.26 |
| Bromobenzene | < 0.24 | 0.075 | 0.24 |
| Bromodichloromethane | < 0.11 | 0.035 | 0.11 |
| n-Butylbenzene | < 0.45 | 0.14 | 0.45 |
| sec-Butylbenzene | < 0.49 | 0.15 | 0.49 |
| tert-Butylbenzene | < 0.4 | 0.12 | 0.4 |
| Carbon Tetrachloride | < 0.5 | 0.16 | 0.5 |
| Chlorobenzene | < 0.27 | 0.086 | 0.27 |
| Chloroethane | < 0.5 | 0.39 | 1.3 |
| Chloroform | < 0.22 | 0.07 | 0.22 |
| Chloromethane | < 1 | 0.88 | 3.1 |
| 2-Chlorotoluene | < 0.65 | 0.21 | 0.65 |
| 4-Chlorotoluene | < 0.19 | 0.06 | 0.19 |
| 1,2-Dibromo-3-Chloropropane | < 1 | 0.83 | 2.7 |
| Dibromochloromethane | < 0.09 | 0.028 | 0.09 |
| 1,2-Dichlorobenzene | < 0.11 | 0.035 | 0.11 |
| 1,3-Dichlorobenzene | < 0.83 | 0.23 | 0.83 |
| 1,4-Dichlorobenzene | < 0.13 | 0.039 | 0.13 |
| Dichlorodifluoromethane | < 5.4 | 1.7 | 5.4 |
| 1,1-Dichloroethene | < 0.37 | 0.12 | 0.37 |
| 1,2-Dichloroethane | < 0.86 | 0.27 | 0.86 |
| 1,1-Dichloroethane | < 0.27 | 0.084 | 0.27 |
| cis 1,2-Dichloroethene | < 0.29 | 0.092 | 0.29 |
| trans-1,2-dichloroethene | < 0.23 | 0.072 | 0.23 |
| 1,2-Dichloropropane | < 0.15 | 0.046 | 0.15 |
| 1,3-DCP, Tetrachloroethene | < 0.56 | 0.17 | 0.56 |

Fluorobenzene Surrogate 108 % Rec.
1,4-Dichlorobutane Surrogate 92 % Rec.
Sample pH 1.6

| ANALYTE | RESULT | MDL UG/L | PQL UG/L |
|---------------------------|--------|-------------|-------------|
| 2,2-Dichloropropane | < 1 | 0.63 | 2.2 |
| Di-isopropyl Ether | < 0.38 | 0.12 | 0.38 |
| Ethylbenzene | < 0.32 | 0.1 | 0.32 |
| EDB (1,2-Dibromoethane) | < 0.08 | 0.025 | 0.08 |
| Hexachlorobutadiene | < 0.35 | 0.11 | 0.35 |
| Isopropylbenzene | < 0.36 | 0.11 | 0.36 |
| p-Isopropyltoluene | < 0.46 | 0.15 | 0.46 |
| Methylene Chloride | < 4 | 0.29 | 0.91 |
| MTBE | < 0.22 | 0.069 | 0.22 |
| Naphthalene | < 0.41 | 0.13 | 0.41 |
| n-Propylbenzene | < 0.41 | 0.13 | 0.41 |
| 1,1,2,2-Tetrachloroethane | < 0.31 | 0.099 | 0.31 |
| Tetrachloroethene | < 0.56 | 0.17 | 0.56 |
| Toluene | < 0.69 | 0.22 | 0.69 |
| 1,2,3-Trichlorobenzene | < 1 | 0.31 | 1.1 |
| 1,2,4-Trichlorobenzene | < 0.91 | 0.26 | 0.91 |
| 1,1,1-Trichloroethane | < 0.63 | 0.2 | 0.63 |
| 1,1,2-Trichloroethane | < 0.17 | 0.055 | 0.17 |
| Trichloroethene | < 0.18 | 0.055 | 0.18 |
| Trichlorofluoromethane | < 1.4 | 1.4 | 4.4 |
| 1,2,4-Trimethylbenzene | < 0.57 | 0.18 | 0.57 |
| 1,3,5-Trimethylbenzene | < 0.57 | 0.18 | 0.57 |
| Vinyl Chloride | < 0.54 | 0.17 | 0.54 |
| m&p-Xylene | < 0.9 | 0.28 | 0.9 |
| o-Xylene | < 0.33 | 0.1 | 0.33 |

MDL = Method Detection Limit

PQL = Practical Quantitation Limit

NA = Not Applicable

Authorized Signature

**Analytical Laboratory**

1090 Kennedy Ave. Kimberly, WI 54136
414-735-8295

WI DNR Certified Lab #445027660

QC Summary**Method 8021 Volatile Organic Compounds**

Project #: CLW131246
Sample ID: Trip

Report Date: 24-Sep-96
Lab Code: 5014513F

| ANALYTE | INITIAL CALIBRATION | KNOWN STANDARD | MATRIX SPIKE | REPLICATE SPIKE | BLANK | PID SURROGATE | HALL SURROGATE |
|-----------------------------|---------------------|----------------|--------------|-----------------|-------|---------------|----------------|
| Benzene | P | P | P | P | P | P | P |
| Bromobenzene | P | P | P | P | P | P | P |
| Bromodichloromethane | P | P | P | F | P | P | P |
| n-Butylbenzene | P | P | P | P | P | P | P |
| sec-Butylbenzene | P | P | P | P | P | P | P |
| tert-Butylbenzene | P | P | P | P | P | P | P |
| Carbon Tetrachloride | P | P | P | P | P | P | P |
| Chlorobenzene | P | P | P | P | P | P | P |
| Chloroethane | P | P | P | P | P | P | P |
| Chloroform | P | P | P | P | P | P | P |
| Chloromethane | P | F | P | P | P | P | P |
| 2-Chlorotoluene | P | P | P | F | P | P | P |
| 4-Chlorotoluene | P | P | P | F | P | P | P |
| 1,2-Dibromo-3-Chloropropane | P | F | P | F | P | P | P |
| Dibromochloromethane | P | P | P | F | P | P | P |
| 1,2-Dichlorobenzene | P | P | P | P | P | P | P |
| 1,3-Dichlorobenzene | P | P | P | P | P | P | P |
| 1,4-Dichlorobenzene | P | P | P | F | P | P | P |
| Dichlorodifluoromethane | P | F | P | P | P | P | P |
| 1,1-Dichloroethane | P | P | P | P | P | P | P |
| 1,2-Dichloroethane | P | P | P | F | P | P | P |
| 1,1-Dichloroethene | P | F | P | P | P | P | P |
| cis-1,2-Dichloroethene | P | P | P | P | P | P | P |
| trans-1,2-Dichloroethene | P | F | P | P | P | P | P |
| 1,2-Dichloropropene | P | P | P | F | P | P | P |
| 1,3-Dichloropropene | P | P | P | F | P | P | P |
| 2,2-Dichloropropene | P | F | P | F | P | P | P |
| Di-Isopropyl Ether | P | P | P | P | P | P | P |
| Ethylbenzene | P | P | P | P | P | P | P |
| EDB (1,2-Dibromoethane) | P | P | P | F | P | P | P |
| Hexachlorobutadiene | P | P | P | F | P | P | P |
| Isopropylbenzene | P | P | P | P | P | P | P |
| p-Isopropyltoluene | P | P | P | P | P | P | P |
| Methylene Chloride | P | F | P | P | P | P | P |
| MTBE | P | P | P | P | P | P | P |
| Naphthalene | P | P | P | P | P | P | P |
| n-Propylbenzene | P | P | P | P | P | P | P |
| 1,1,2,2-Tetrachloroethane | P | P | P | F | P | P | P |
| Tetrachloroethene | P | P | P | P | P | P | P |
| Toluene | P | P | P | P | P | P | P |
| 1,2,3-Trichlorobenzene | P | P | P | F | P | P | P |
| 1,2,4-Trichlorobenzene | P | P | P | F | P | P | P |
| 1,1,1-Trichloroethane | P | P | P | P | P | P | P |
| 1,1,2-Trichloroethane | P | P | P | F | P | P | P |
| Trichloroethene | P | P | P | P | P | P | P |
| Trichlorofluoromethane | P | F | F | P | P | P | P |
| 1,2,4-Trimethylbenzene | P | P | P | P | P | P | P |
| 1,3,5-Trimethylbenzene | P | P | P | P | P | P | P |
| Vinyl Chloride | P | F | P | P | P | P | P |
| m & p-Xylene | P | P | P | P | P | P | P |
| o-Xylene | P | P | P | P | P | P | P |

P = Passed QC limits.

F = Failed QC limits.

NA = Not Applicable

VOC analysis detected unidentified peaks.

Authorized Signature

**Analytical Laboratory**

1090 Kennedy Ave. Kimberly, WI 54136
414-735-8295

WI DNR Certified Lab #445027660

Method 8021 Volatile Organic Compounds

BRIAN KRANZ
NORTHERN ENVIRONMENTAL
1214 W VENTURE COURT
MEQUON WI 53092

Report Date: 24-Sep-96
Analyzed By: R. Everson

Project #: CLW131246
Project : Cedarburg
Sample ID: FB-1246
Lab Code: 5014513G
Sample Type: Water
Sample Date: 13-Sep-96
Date Analyzed: 19-Sep-96

| ANALYTE | RESULT | MDL UG/L | PQL UG/L |
|-----------------------------|--------|-------------|-------------|
| Benzene | < 0.26 | 0.082 | 0.26 |
| Bromobenzene | < 0.24 | 0.075 | 0.24 |
| Bromodichloromethane | < 0.11 | 0.035 | 0.11 |
| n-Butylbenzene | < 0.45 | 0.14 | 0.45 |
| sec-Butylbenzene | < 0.49 | 0.15 | 0.49 |
| tert-Butylbenzene | < 0.4 | 0.12 | 0.4 |
| Carbon Tetrachloride | < 0.5 | 0.16 | 0.5 |
| Chlorobenzene | < 0.27 | 0.086 | 0.27 |
| Chloroethane | < 0.5 | 0.39 | 1.3 |
| Chloroform | < 0.22 | 0.07 | 0.22 |
| Chloromethane | < 1 | 0.88 | 3.1 |
| 2-Chlorotoluene | < 0.65 | 0.21 | 0.65 |
| 4-Chlorotoluene | < 0.19 | 0.06 | 0.19 |
| 1,2-Dibromo-3-Chloropropane | < 1 | 0.83 | 2.7 |
| Dibromochloromethane | < 0.09 | 0.028 | 0.09 |
| 1,2-Dichlorobenzene | < 0.11 | 0.035 | 0.11 |
| 1,3-Dichlorobenzene | < 0.83 | 0.23 | 0.83 |
| 1,4-Dichlorobenzene | < 0.13 | 0.039 | 0.13 |
| Dichlorodifluoromethane | < 5.4 | 1.7 | 5.4 |
| 1,1-Dichloroethene | < 0.37 | 0.12 | 0.37 |
| 1,2-Dichloroethane | < 0.86 | 0.27 | 0.86 |
| 1,1-Dichloroethane | < 0.27 | 0.084 | 0.27 |
| cis 1,2-Dichloroethene | < 0.29 | 0.092 | 0.29 |
| trans-1,2-dichloroethene | < 0.23 | 0.072 | 0.23 |
| 1,2-Dichloropropane | < 0.15 | 0.046 | 0.15 |
| 1,3-DCP, Tetrachloroethene | < 0.56 | 0.17 | 0.56 |

| ANALYTE | RESULT | MDL UG/L | PQL UG/L |
|---------------------------|--------|-------------|-------------|
| 2,2-Dichloropropane | < 1 | 0.63 | 2.2 |
| Di-isopropyl Ether | < 0.38 | 0.12 | 0.38 |
| Ethylbenzene | < 0.32 | 0.1 | 0.32 |
| EDB (1,2-Dibromoethane) | < 0.08 | 0.025 | 0.08 |
| Hexachlorobutadiene | < 0.35 | 0.11 | 0.35 |
| Isopropylbenzene | < 0.36 | 0.11 | 0.36 |
| p-Isopropyltoluene | < 0.46 | 0.15 | 0.46 |
| Methylene Chloride | < 4 | 0.29 | 0.91 |
| MTBE | < 0.22 | 0.069 | 0.22 |
| Naphthalene | < 0.41 | 0.13 | 0.41 |
| n-Propylbenzene | < 0.41 | 0.13 | 0.41 |
| 1,1,2,2-Tetrachloroethane | < 0.31 | 0.099 | 0.31 |
| Tetrachloroethene | < 0.56 | 0.17 | 0.56 |
| Toluene | < 0.69 | 0.22 | 0.69 |
| 1,2,3-Trichlorobenzene | < 1 | 0.31 | 1.1 |
| 1,2,4-Trichlorobenzene | < 0.91 | 0.26 | 0.91 |
| 1,1,1-Trichloroethane | < 0.63 | 0.2 | 0.63 |
| 1,1,2-Trichloroethane | < 0.17 | 0.055 | 0.17 |
| Trichloroethene | < 0.18 | 0.055 | 0.18 |
| Trichlorofluoromethane | < 1.4 | 1.4 | 4.4 |
| 1,2,4-Trimethylbenzene | < 0.57 | 0.18 | 0.57 |
| 1,3,5-Trimethylbenzene | < 0.57 | 0.18 | 0.57 |
| Vinyl Chloride | < 0.54 | 0.17 | 0.54 |
| m&p-Xylene | < 0.9 | 0.28 | 0.9 |
| o-Xylene | < 0.33 | 0.1 | 0.33 |

MDL = Method Detection Limit

PQL = Practical Quantitation Limit

NA = Not Applicable

Fluorobenzene Surrogate 108 % Rec.
1,4-Dichlorobutane Surrogate 89 % Rec.
Sample pH 1.6

Authorized Signature

**Analytical Laboratory**

1090 Kennedy Ave. Kimberly, WI 54136
414-735-8295

WI DNR Certified Lab #445027660

QC Summary**Method 8021 Volatile Organic Compounds**

Project #: CLW131246 Report Date: 24-Sep-96
Sample ID: FB-1246 Lab Code: 5014513G

| ANALYTE | INITIAL CALIBRATION | KNOWN STANDARD | MATRIX SPIKE | REPLICATE SPIKE | BLANK | PID SURROGATE | HALL SURROGATE |
|-----------------------------|---------------------|----------------|--------------|-----------------|-------|---------------|----------------|
| Benzene | P | P | P | P | P | P | P |
| Bromobenzene | P | P | P | P | P | P | P |
| Bromodichloromethane | P | P | P | F | P | P | P |
| n-Butylbenzene | P | P | P | P | P | P | P |
| sec-Butylbenzene | P | P | P | P | P | P | P |
| tert-Butylbenzene | P | P | P | P | P | P | P |
| Carbon Tetrachloride | P | P | P | P | P | P | P |
| Chlorobenzene | P | P | P | P | P | P | P |
| Chloroethane | P | P | P | P | P | P | P |
| Chloroform | P | P | P | P | P | P | P |
| Chloromethane | P | F | P | P | P | P | P |
| 2-Chlorotoluene | P | P | P | F | P | P | P |
| 4-Chlorotoluene | P | P | P | F | P | P | P |
| 1,2-Dibromo-3-Chloropropane | P | F | P | F | P | P | P |
| Dibromochloromethane | P | P | P | F | P | P | P |
| 1,2-Dichlorobenzene | P | P | P | P | P | P | P |
| 1,3-Dichlorobenzene | P | P | P | P | P | P | P |
| 1,4-Dichlorobenzene | P | P | P | F | P | P | P |
| Dichlorodifluoromethane | P | F | P | P | P | P | P |
| 1,1-Dichloroethane | P | P | P | F | P | P | P |
| 1,2-Dichloroethane | P | P | P | F | P | P | P |
| 1,1-Dichloroethene | P | F | P | P | P | P | P |
| cis-1,2-Dichloroethene | P | P | P | P | P | P | P |
| trans-1,2-Dichloroethene | P | F | P | P | P | P | P |
| 1,2-Dichloropropane | P | P | P | F | P | P | P |
| 1,3-Dichloropropane | P | P | P | F | P | P | P |
| 2,2-Dichloropropane | P | F | P | F | P | P | P |
| Di-Isopropyl Ether | P | P | P | P | P | P | P |
| Ethylbenzene | P | P | P | P | P | P | P |
| EDB (1,2-Dibromoethane) | P | P | P | F | P | P | P |
| Hexachlorobutadiene | P | P | P | F | P | P | P |
| Isopropylbenzene | P | P | P | P | P | P | P |
| p-Isopropyltoluene | P | P | P | P | P | P | P |
| Methylene Chloride | P | F | P | P | P | P | P |
| MTBE | P | P | P | P | P | P | P |
| Naphthalene | P | P | P | P | P | P | P |
| n-Propylbenzene | P | P | P | P | P | P | P |
| 1,1,2,2-Tetrachloroethane | P | P | P | F | P | P | P |
| Tetrachloroethene | P | P | P | P | P | P | P |
| Toluene | P | P | P | P | P | P | P |
| 1,2,3-Trichlorobenzene | P | P | P | F | P | P | P |
| 1,2,4-Trichlorobenzene | P | P | P | F | P | P | P |
| 1,1,1-Trichloroethane | P | P | P | P | P | P | P |
| 1,1,2-Trichloroethane | P | P | P | F | P | P | P |
| Trichloroethene | P | P | P | P | P | P | P |
| Trichlorofluoromethane | P | F | F | P | P | P | P |
| 1,2,4-Trimethylbenzene | P | P | P | P | P | P | P |
| 1,3,5-Trimethylbenzene | P | P | P | P | P | P | P |
| Vinyl Chloride | P | F | P | P | P | P | P |
| m & p-Xylene | P | P | P | P | P | P | P |
| o-Xylene | P | P | P | P | P | P | P |

P = Passed QC limits.

F = Failed QC limits.

NA = Not Applicable

VOC analysis detected unidentified peaks.

Authorized Signature

