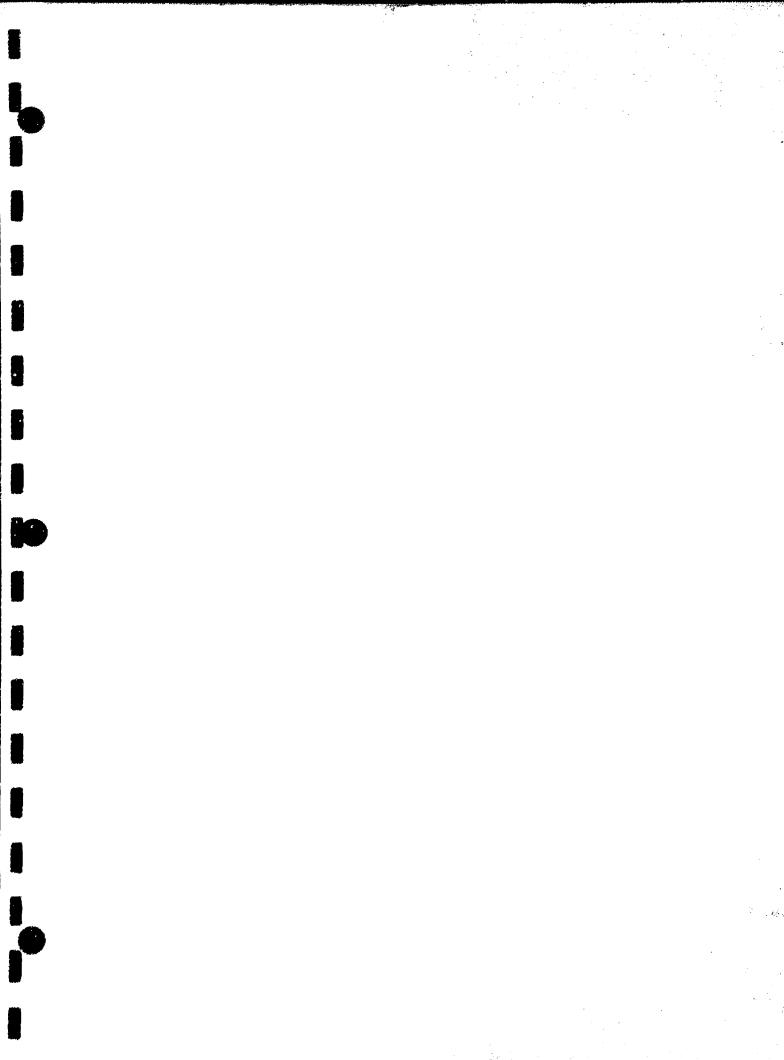
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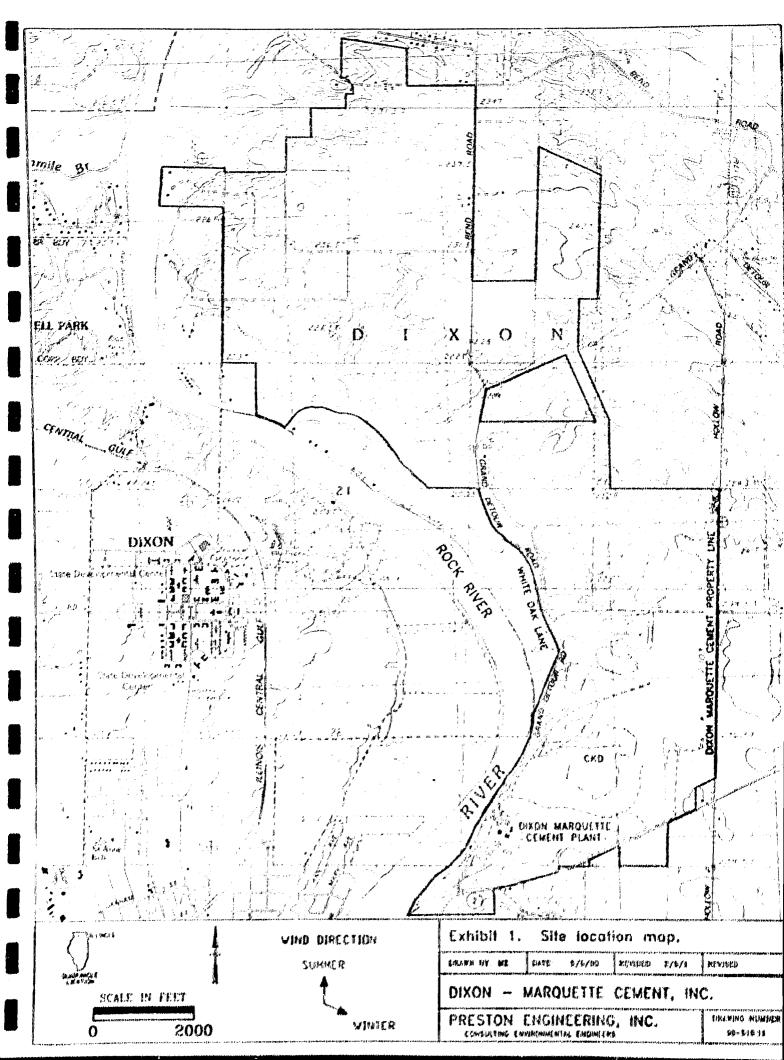
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March 15, 2001

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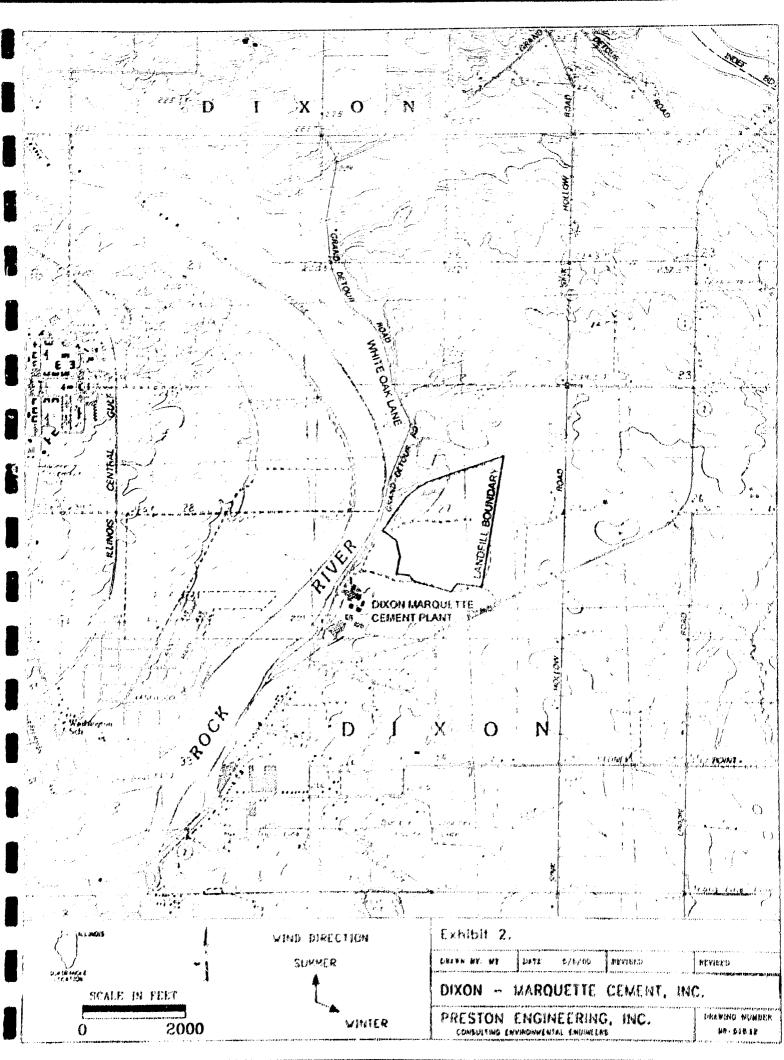
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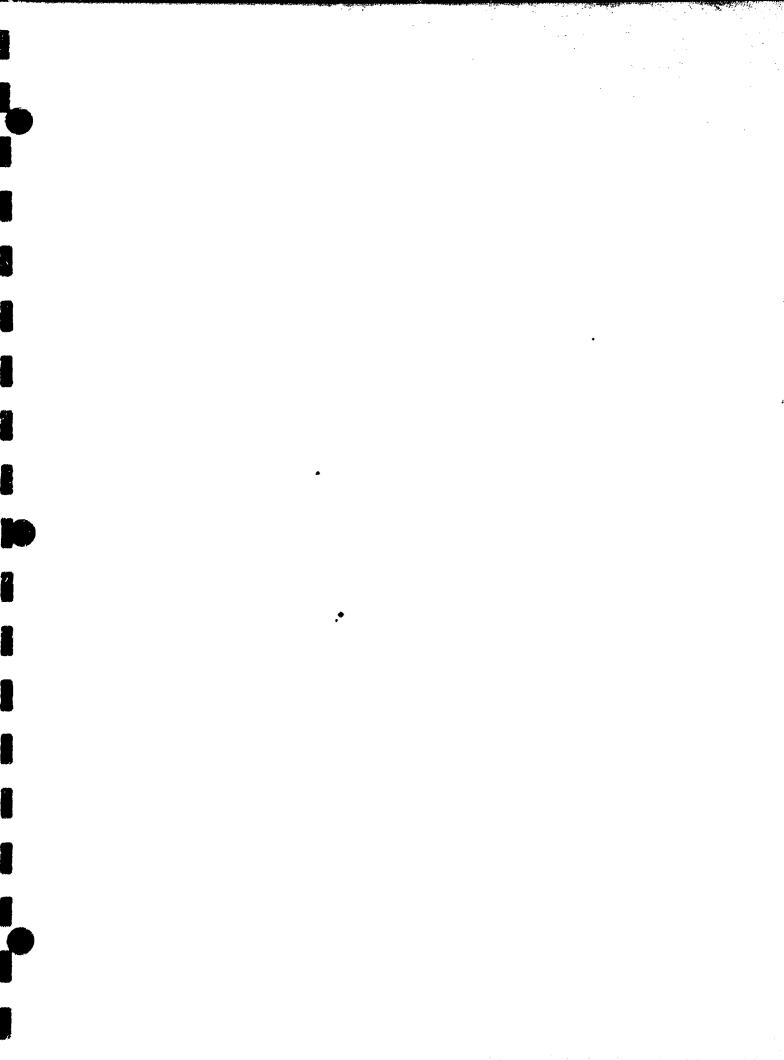


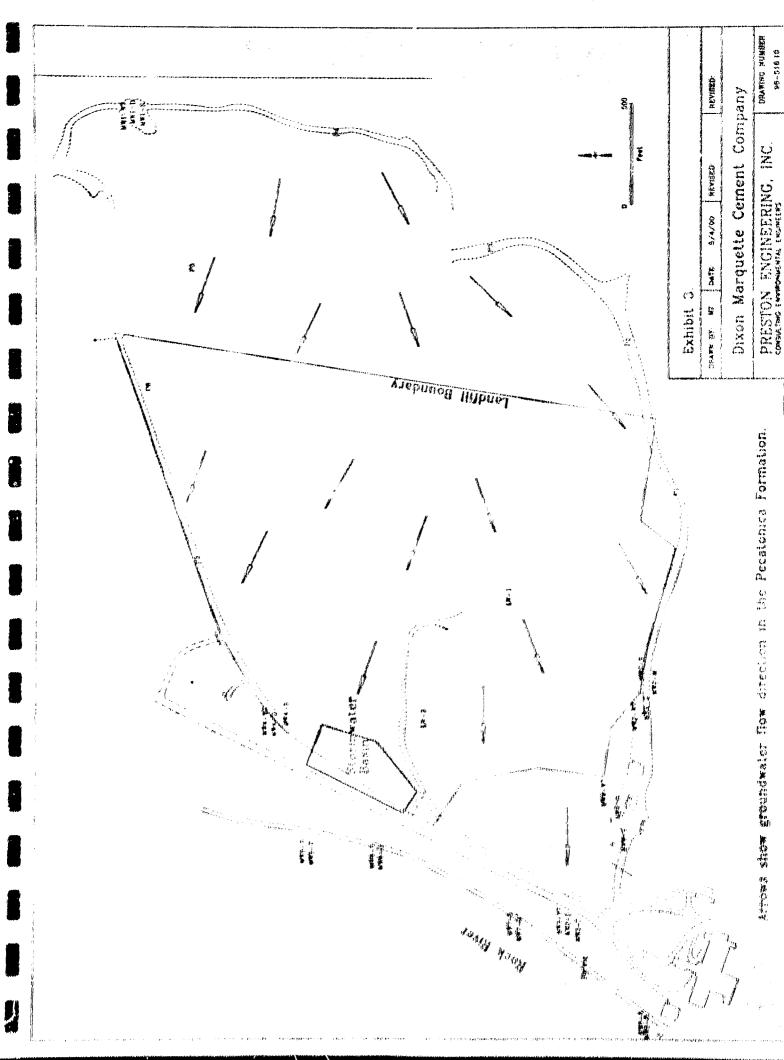


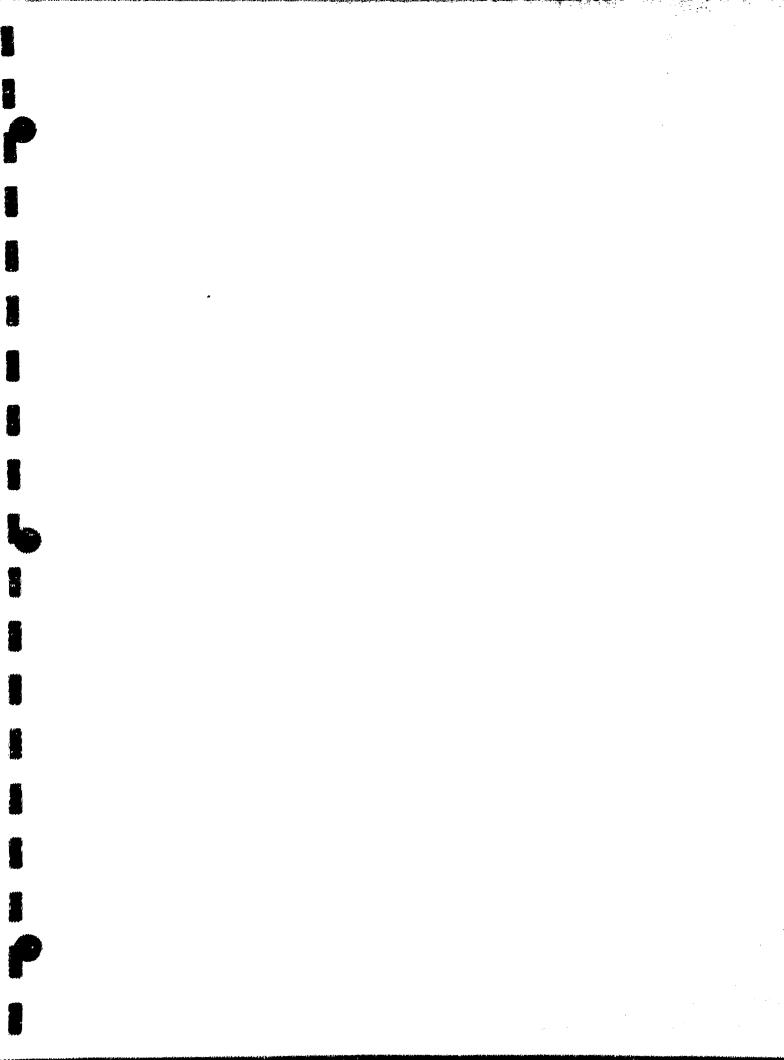
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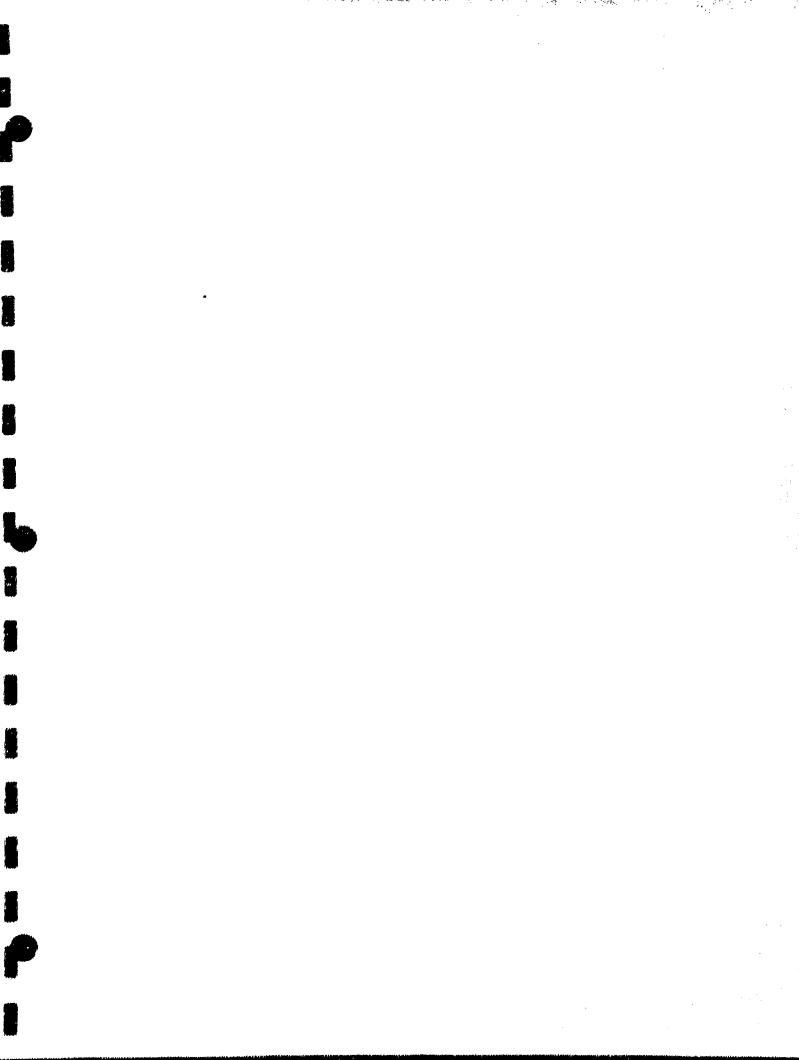
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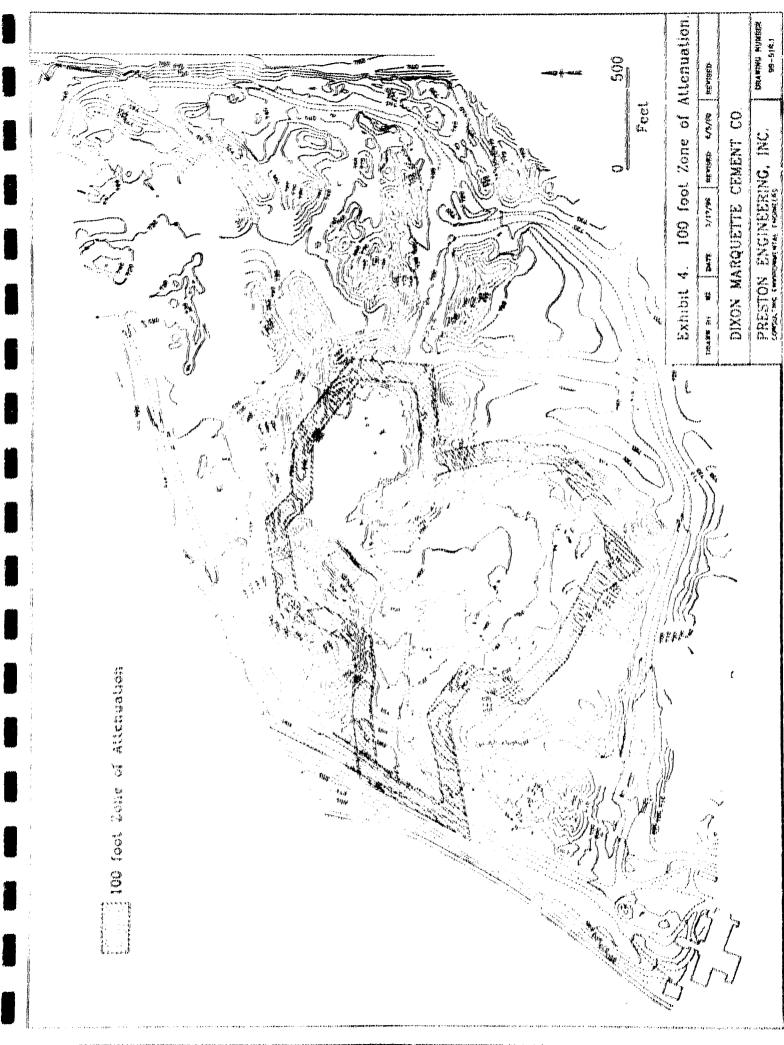


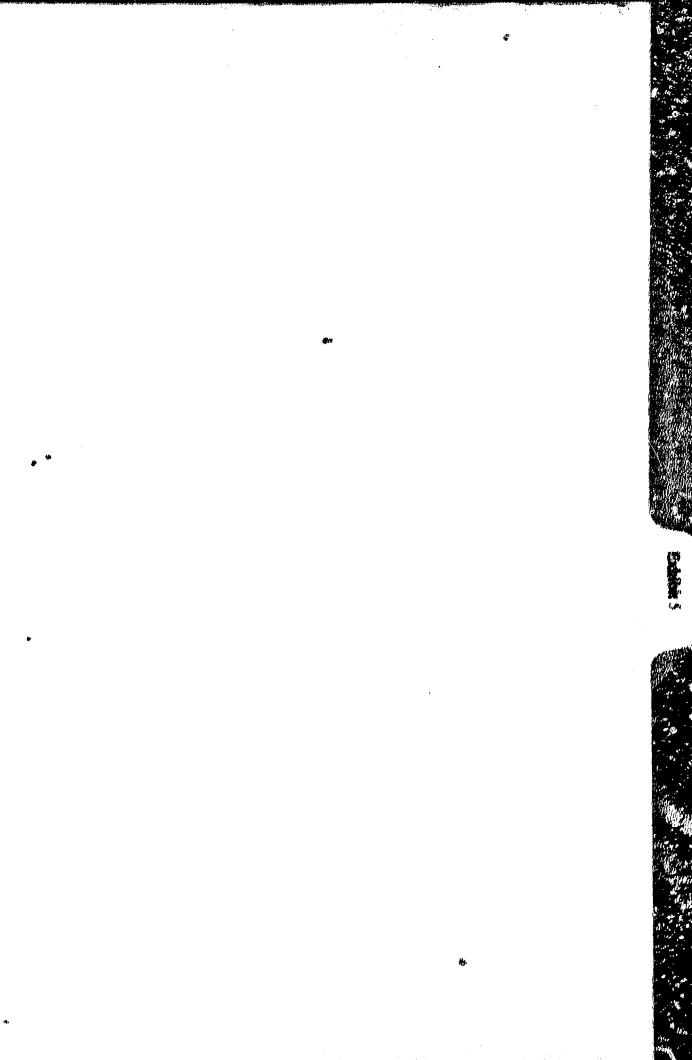












# Hydrogeologic Report Dixon Marquette Cement March 2001

Prepared by

Presson Engineering, Inc.
4436 N. Brady Street, Davenport, Iowa

Mark W. Zell

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#### 1.0 Introduction

This report describes the geologic and hydrologic conditions at the Dixon Marquette Cement Company site near Dixon, Illinois, and the surrounding area. Information from published material and from field studies at the site that were completed by Preston Engineering, Inc. and Versar, Inc. was used. Wells and piezometers were installed in and around the landfill. The wells and the Rock Biwer were sampled to establish the chemical characteristics of the groundwater and surface water. Hydraulic conductivity tests were conducted on the monitoring wells. A study of bedrock structure and joint patterns was completed. The information from the published sources and from the site investigation is presented and interpreted in this report.

#### 2.0 Geologie Setting

#### 2.1 Regional Geology

The area within 10 miles of Dixon has Ordovreian age bedrock near the surface, with exposures along the Rock River. The bedrock includes (in order of increasing age) the Maquoketa, Galena, Platteville, Ameell, and Prairie du Chien Groups, and is mossy limestone, dolomite, sandstone, and nome shale. The Maquoketa and Galena Groups consist of shale, limestone, and dolomite, and they outerop to the southwest of Dixon. The Platteville Group consists of limestone and dolomite, and it materials in the Dixon area. It is primarily limestone of the Mifflin Formation in the Platteville Group that is mined at the Dixon Marquette Cement plant. To the northeast of Dixon, sandstone of the Ancell Group and dolomite of the Prairie du Chien Group outerop. Lonally, Pennsylvanian shale occurs in fissures and sinkholes in the Platteville Group.

The acqueence of rocks chooses increasing age from somthwest to northeast, and the rocks dip to the sandhwest. The nocks lie on the flank of the Ashnon Arch and at the continuence with of the Wisconsin Arch. The nocks dip areas from the center of the Ashnon Arch, which trends northwest to sendium to the name and of Discon. Panisheling the nich is the Sandarch fault zone, which runs from near Despots to seem Marchanian. On the application of all some wide of the fault. Caption rocks have been bronged to the southern. The total range is 0.5 to 2 wides wide and is about 85 miles have been bronged to the surface displacement along the range is the about 800 feet.

The Rock River flows to the southwest near Dixon, and it has cut into the bedrock. The St. Peter sandstone is exposed along the river near Grand Detour, and the Platteville Group is exposed near Dixon. The river makes a sharp horseshoe-shaped meander near Grand Detour, possibly as a consequence of joint systems in the bedrock.

Unconsolidated Holocene deposits consist of alluvium in the Rock River valley, glacial till deposits, and wind-blown loess.

#### 2.2 Site Geology

The surface material has been disturbed in and near the CKD disposal area. Overlying the bedrock are two types of material: soil and rock spoil from mining operations, typically placed in windrows on top of the Pecatonica Formation, and cement kiln dust, clinker, and other debris, typically placed on top of the Pecatonica Formation and the spoil. The thickness of the spoil reaches 80 to 90 feet in the western part of the landfill. The thickness of the CKD reaches about 60 feet near the center of the landfill.

A description of the stratigraphic column at the Dixon Marquette Cement plant was published in the Illinois State Geological Survey Circular 502 (Willman and Kolata, 1978). In the circular, the Grand Detour, Mifflin, and Pecatonica Formations are described as they appear in outcrop. These three formations are exposed at the site, and are described below based on descriptions by Willman and Kolata, field observations of outcrops by Preston Engineering, Inc., and examination of rock cores from holes drilled for monitoring wells.

The Grand Detour Formation is approximately 32 feet thick, and consists of the Forceston member (12 feet thick) and the Stillman member (20 feet thick). The Forceston member is a gray argillaceous fossiliferous limestone, and the Stillman member is a gray limestone wan dolomite mostling. There are shaly partings in both members. The Grand Detour Formation is exposed in elitis on the west side of Sink Hollow Road where it and the underlying Millin Formation have been mined.

The Mifflin Formation is the primary source of the limestone used by Dixon Marquetic in the manufacture of Portland cement. It consists of about 18 feet of gray and yellow argillaceous fossiliferous limestone with shale partings. There is a 4 to 6 inch thick layer of dolomitic shale at the top, underlain by a layer of bentonite that is usually less than 0.5 inch thick. The Mifflin Formation has thin wavy bedding, usually 0.5 to 3 inches thick, and has numerous vertical joi is. The limestone within the Mifflin Formation typically has five percent or less magne um carbonate. It is exposed on the west side of Sink Hollow Road, on the east side of White Oak Lane, and in outliers in the mined areas around the landfill where the rock was not suitable for use in cement manufacturing or not accessible.

The Pecatonica Formation consists of about 34 feet of limestone and dolomite. Four members are present at the site: the Medusa member, a dolomite-mottled limestone with a corrosion surface at the top; the New Glarus member, consisting of dolomite and dolomite-mottled limestone with a thin shaly bed at the top; the Dane member, a limestone; and the Chana member, a sandy dolomite. The bedding is thicker than in the Mifflin I Immation, and is about 2 to 12 inches thick. It has wavy shaly bedding in places, and some stylolites, suggesting that some carbonate has been dissolved. The lower Pecatonica is vuggy in places, which increases the porosity of the rock. The dolomite in the Pecatonica Formation typically has between 5 and 25 percent magnesium carbonate. The top of the Pecatonica is exposed in much of the area around the CKD landfill at Dixon Marquette, and probably directly underlies most of the landfill. It is also exposed along the Rock River west of the landfill. The Pecatonica was exposed around the landfill as a result of mining of the overlying Mifflin Formation. Dixon Marquette uses the Pecatonica Formation as a source of construction aggregate.

The Platteville Group rocks have numerous joints, and some can be traced on horizontal surfaces for several hundred feet. The joints are enlarged in places, which provide local areas of extremely high permeability. The dominant bearing of the joints is between N80°E and N80°W. There is a secondary trend with bearings between N10°W and N10°E. Over 90 percent of the joints are within 5° of vertical.

in most areas in and immediately around the landfill the rock overlying the Pecatonica Formation has been mined. The mined rock was blasted prior to removal. The blasting procedure consisted of drilling vertical holes from the surface to the top of the Pecatonica Formation, and detonating explosives in the holes. The holes were horizontally spaced about 10 feet apart. The rock overlying the Pecatonica was fractured by the blasting and was hauled to the plant. The blasting was designed to allow removal of the Mifflin Formation without significant damage to the Pecatonica Formation. The top of the Pecatonica Formation has been used as a work surface by heavy equipment to load and haul the blasted rock, and it remains largely intact today. The effects of blasting can be seen in the Pecatonica Formation, however. Numerous circular potholes, roughly 3 to 4 feet in diameter, mark the top of the rock. The holes are usually no deeper than about 0.5 foot, and are spaced about 10 feet apart. Fractures radiate in a spoke-like pattern from many of the holes, and intersect fractures from neighboring holes. The potholes originated at the base of blast holes made in the overlying formations, and the fractures resulted from the blasting. The vertical extent of the blast fractures was not observed.

The removal of the Mifflin Formation in the mined areas has revealed the structure or, top of the Pecatonica Formation. The Pecatonica Formation dips gently westward, but locally there are high spots or small domes. The domes range from about 50 fect to several hundred feet in diameter, and have up to about 10 feet of relief in the larger domes. Typically the Mifflin Formation above the domes is extensively jointed and has a chemical composition that makes it unsuitable for cement production.

Underlying the Pecatonica and not exposed at the site is the Ancell Group, which consists of the Glenwood and St. Peter Formations. The Glenwood Formation consists of 8 to 10 feet of gray fine quartz sandstone and gray siltstone. The sandstone is mottled and well cemented. The quartz grains are poorly sorted and range in size from 0.1 mm to 0.5 mm. The poor sorting distinguishes the sandstone from the underlying St. Peter Formation, which is well sorted and has sand grains of about 0.5 mm. Within the Glenwood Formation sandstone are two siltstone layers that are 1 to 2 feet thick. The upper siltstone layer is sandy at the top. The siltstone separates easily along bedding planes, and is more friable than the sandstone. Outcrops of the Glenwood in the area are rare, propably because of its thickness and because the siltstone is easily eroded. The elevation of

 the top of the Glenwood Formation was determined at four monitoring well locations around the landfill, and it was found that the rock strikes N 5°W and has a dip of 0.9° to the west.

The §!. Peter Formation is a uniform nearly pure quartz sandstone. A water supply well drilled at the Dixon Marquette site in 1992 encountered 359 feet of sandstone of the St. Peter Formation. It consists almost entirely of fine to medium, well sorted, well rounded, poorly cemented quartz grains. There are some shaly layers and carbonate-cemented layers near the top of the formation.

Cross sections of the site are included in this report. A description of geologic materials at the monitoring well locations is also included.

#### 3.0 Hydrologic Setting

#### 3.1 Regional Hydrology

Surface water drainage in the area around Dixon is dominated by the Rock River and its tributaries. The Rock River in Illinois flows to the southwest from north of Rockford to Rock Island. Tributaries flowing from the upland areas create a dendritic drainage pattern. There is a maximum of about 190 feet of relief within 3 miles of downtown Dixon.

The important aquifers in the area are the St. Peter Formation and the Mt. Simon Formation, both of which are sandstone. The St. Peter Formation is exposed to the northeast of Dixon, in the area around Grand Detour. The exposed area is probably a recharge zone for the aquifer. The rock dips gently to the southwest. The Mt. Simon Formation consists of about 1500 feet of sandstone with some shale, with the top of the formation found at about 1000 feet deep in the Dixon area. The City of Dixon maintains several water supply wells one to two miles from the landfill. They have a total depth of 1260 to 1870 feet, and obtain water from the St. Peter, Eminence, and Mt. Simon Formations.

#### 3.2 Site Hydrology

Prior to mining, the area of the landfill was hilly with a northwest-trending valley leading to the Ruck River. There was 80 to 100 feet of relief. The extreme west side of the landfill was used for

railroad sidings. The Mifflin Formation may have been thin or absent on the former valley floor on the northwest side of the landfill, and possibly no rock was mined in that area. Mining probably progressed up the dip of the rocks to the east and northeast to aid in draining the work area.

Currently, surficial drainage consists of small streams that parallel roads around the landfill. The configuration of mined and filled areas around the landfill prevents good drainage, and small ponds have formed. On the east side of the landfill groundwater exits the rock faces near Sink Hollow Road and forms ponds. Some of the ponds have outlets to ephemeral streams, and others do not appear to have an outlet. Those with no outlet must drain through evaporation and infiltration. In 1999, a storm water basin was constructed on the west side of the landfill. The basin was designed to restrict storm water runoff from flowing directly into the Rock River and decrease the amount of solids being discharged.

The Pecatonica Formation is exposed at the surface or underlies most of the fill material in the landfill. The rock is extensively jointed, and surface water and groundwater can seep into the rock via the joints. The vertical extent of the joints was not determined during the field investigation. On the east side of the landfill along a road, water was observed to be flowing into an opening in the rock. The flow rate into the opening during a period of rain in May 1998 was estimated at 10 to 20 gallons per minute. The water entering the hole was turbid and pale yellow.

The Pecatonica is exposed along the Rock River to the west of the landfill. Springs occur along the exposures, and the water flows into the river. The largest of the springs in this area, located along White Oak Lane east-southeast of the landfill, exits a joint in the rock in the upper part of the Pecatonica Formation, about 5 feet below the contact with the overlying Mifflin Formation. The flow during several days in March 1998 was estimated at 100 gallons per minute, or about 160 acre-feet per year. The flow approximately doubled shortly after a period of heavy rain in May 1998. The water flowing from the spring was observed to be clear on several occasions in March, and was turbid and pale yellow on two occasions in May. Because of the fluctuations in discharge and changes in turbidity that are apparently associated with precipitation, it is thought that there exist connections between surface water and the joint from which the spring issues.

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In some places the Pecatonica Formation contains shaly or clay layers that impede the vertical migration of groundwater. Groundwater was observed to be discharging immediately above a thin clay layer in the Pecatonica Formation on the east side of the coal storage area north of the landfill. This clay layer was similar in appearance and thickness to the bentonite layer that occurs near the top of the Mifflin Formation. Groundwater was also observed to be exiting the rock above the bentonite layer in the Mifflin Formation near Sink Hollow Road. The groundwater discharges at these points indicates that layers of clay less than 0.5 inch thick can be effective in redirecting the vertical flow of groundwater in the Platteville Group.

The uppermost aquifer at the site is almost entirely contained within the Pecatonica Formation. The top of the aquifer is most commonly about one to five feet below the top of the Pecatonica Formation, or what remains of the formation in areas where the top portion of it was removed during mining. In areas with thick CKD deposits the top of the aquifer is about five feet above the Pecatonica Formation, and is within the CKD. The bottom of the aquifer coincides with the contact between the Pecatonica Formation and the underlying Glenwood Formation. The depth of this contact ranges from about 34 to 100 feet. The Glenwood Formation consists of sandstone and shale, and it acts as an aquitard. The depth of the top of the aquifer ranges from about zero to 90 feet at the facility. The thickness of the aquifer ranges from about 30 to 43 feet, and is thickest where it underlies the CKD disposal area.

Quarterly measurements of the water table elevation in the uppermost aquifer at four monitoring well locations around the CKD disposal area were made from April 1998 to November 2000. The measurements indicated that the water table had a gradient of about 0.02, and was flowing west and southwest toward the Rock River. The difference between the highest and lowest measured elevation of the water table at the casternmost well, MW1-S, was 0.13 feet. At the westernmost well, MW3-S, the difference was 2.35 feet. The potentiometric surface map included in this report shows a slight mound of water beneath the CKD landfill. If groundwater is mounded in the CKD area, the flow direction should be slightly northwest and southwest, as well as west, from the CKD area. The water table and the direction of groundwater flow in the uppermost aquifer appear to be stable throughout the year, based on the quarterly measurements. Short term temporal variations in the elevation of the water table have not been investigated.

Quarterly measurements of the potentiometric surface were also made in wells screened in the St. Peter Sandstone. Data from the four deep wells indicated that the water table has a westward gradient. The most fluctuation over one year of measurements was at well MW3-D, located on the west side of the landfill. The fluctuation was 2.33 feet. The water table and the direction of groundwater flow in the St. Peter aquifer appear to be stable throughout the year, based on the quarterly measurements. Short term temporal variations in the elevation of the potentiometric surface have not been investigated.

The potentiometric surface in both deep and shallow wells apparently dropped by about 1 to 3 feet between Spring 1998 and Summer 1998. The water levels had not recovered to Spring 1998 conditions by February 1999. It is not known if the fluctuation is correlated to a seasonal effect such as water uptake by vegetation or wet and dry periods.

The hydraulic conductivity of the area around the well screens at 22 on site monitoring wells was measured by means of slug tests. Tests were conducted by positioning a pressure transducer beneath the water in the well, quickly submersing a solid slug, and recording water level changes in response. After the water level stopped falling and was at or near the initial static water level, the slug was quickly removed and the response of the water level was recorded again. An In-Situ pressure transducer and data logger were used to record the test data. The data was evaluated using the method of Bouwer and Rice. The data for the test at well MW5-8 did not have sufficient resolution to conduct an adequate analysis. The following table shows the calculated hydraulic conductivity at each of the other wells.

Table 1. Hydraulic Conductivity at Dixon Marquette Cement Company Monitoring Wells

Well	K (feet/minute)	K (centimeters/second)
MW1-S	0.00735	0.0037
MW2-S	0.0000228	0.000012
MW3-S	0.000155	0.000079
MW4-S	0.00195	0.00099
MW6-S	0.0089	0.0045
MW7-S	0.015	0.0078
MW8-S	0.014	0.0073
MW9-S	0.000045	0.000023
MWI-D	0.00024	0.00012
MW2-D	0.0108	0.0055
MW3-D	0.00061	0.00031
MW4-D	0.000817	0.00042
MW5-D	0.031	0.016
MW6-D	0.0061	0.0031
MW7-D	0.0030	0.0015
MW8-D	0.0024	0.0012
MW9-D	0.0028	0.0014
MW2-M	0.000000833	0.00000042
MW2-WT	0.0037	0.0019
MW3-WT	0.00013	0.00066
MW9-WT	0.0089	0.0045

The permeability of the Pecatonica Formation varies greatly both horizontally and vertically, and is probably dependent on the proximity of joints. Tests on monitoring wells installed at the base of the Pecatonica Formation showed that the highest measured hydraulic conductivity was about 300 times more than the lowest conductivity.

Based on the measured hydraulic conductivity and gradient, and estimates of the porosity of the rock, the groundwater flow velocity in the Pecatonica Formation ranges from 0.001 to 38 feet per day. Actual flow rates may be several orders of magnitude higher in areas with large fractures.

The St. Peter Formation underlies the Glenwood Formation, and consists of uniform fine to medium quartz sandstone. It is an important aquifer in the area. Tests on monitoring wells completed in the formation on site (the deep wells) show that the conductivity of the sandstone varies from 0.00024 to 0.031 feet per minute. The water in the St. Peter Formation may be artesian on the west side of the disposal area.

Between the Pecatonica Formation and the St. Peter Formation is sandstone and siltstone of the Glenwood Formation. Injection tests were conducted in the formation at the time of monitoring well installation to determine the permeability of the shale within the Glenwood Formation. The test apparatus consisted of a double packer that was used to isolate the interval of the formation to be tested, a water pump capable of increasing the water pressure between the packers, a pressure gauge that measured the water pressure between the packers, and a flow meter that measured the amount of water moving into the formation. After sealing the packers, the water pressure was increased in increments and then held steady for 30 minutes. At well MW2-D some flow was measured during the test, but it could not be determined if both packers had properly scaled against the borehole wall. Because the bottom packer may have leaked and the water flowed into a more permeable underlying formation, the validity of the injection test at well MW2-i) is in doubt. The test apparatus was modified for tests on other wells by removing the bottom packer, and drilling only to the bottom of the section of rock to be tested. No leakage was seen around the single packer during these tests, and any flow should have been into the rock below the packer. Because no flow was measured during the tests, the calculated conductivities represent an estimated maximum conductivity based on the sensitivity of the flow meter that was used and the pressure that was applied. The results of aquifer tests on the Glenwood Formation are summarized in the following table.

\* Table 2. Hydraulic Conductivity of the Glenwood Formation\*

Well	Hydraulic Conductivity (feet/minute)	Hydraulic Conductivity (cm/second)
MW1-D	<5.83 x 10 <sup>-7</sup>	<2.96 x 10 <sup>-7</sup>
MW2-M	8.33 x 10 <sup>-7</sup>	4.2 x 10 <sup>-7</sup>
MW3-D	<4.12 x 10 <sup>-7</sup>	<2.09 x 10 <sup>-7</sup>
MW4-D	<5.27 x 10 <sup>-7</sup>	<2.67 x 10 <sup>-7</sup>
MW7-D	1.3 x 10 <sup>-4</sup>	6.68 x 10 <sup>-5</sup>
MW8-S	4.4 x 10 <sup>-7</sup>	2.22 x 10 <sup>-7</sup>
MW9	1.4 x 10 <sup>-7</sup>	6.86 x 10 <sup>-7</sup>

Because no flow into the rock was measured during injection tests at walls MWI-D, MWJ-D, and MW4-D, the resulting hydraulic conductivities represent a maximum value based on the sensitivity of the flow measuring device. The hydraulic conductivity at well MW2-M was measured with a slug test. The hydraulic conductivity at wells MW7-D, MW8-S, and MW9 was determined by laboratory analysis of rock core samples.

No evidence of vertical fractures was seen in core samples from the formation. A monitoring well was installed in the Glenwood Formation (MW2-M), and a slug test was conducted. The hydraulic conductivity was calculated to be  $8.3 \times 10^{-7}$  feet per minute, or about 300 to 9000 times less than the overlying dolomite.

There are additional sandstone formations underlying the St. Peter Formation that werve as aquifers, but because of their depth and isolation from the surficial activities at the landfill, they are not discussed here.

Maps of the site showing the potentiometric surface of the aquifers in the Pecatonica and St. Peter Formations are included in this report.

#### 3.3 Climatic Conditions

Information on the climate at Dixon was obtained from the Midwest Climate Center in Springfield, Illinois. Between 1961 and 1990 the average annual temperature was 48.1°F, the

average annual maximum was 58.7°, and the average annual minimum was 37.5°. Extreme temperatures from 1893 to 1996 were -27° and 110°. The average annual precipitation between 1961 and 1990 was 35.58 inches, and between 1893 and 1996 it ranged from 18.49 inches in 1994 to 47.53 inches in 1965. Ten-year 24-hour rainfall events produce between 4.5 and 5.0 inches of rain. The number of days per year with a base temperature of 32° or more averaged 168. Wind speed and direction for Rockford, Illinois (30 miles northeast of Dixon) averaged 9.8 miles per hour, and the average direction was from the west-northwest during the winter and from the south at other times.

#### 4.0 Monitoring Well Installation

During March and April 1998, nine monitoring wells were installed around the CKD disposal site. The wells were installed in pairs at three sites, and three wells were installed at one site. Additional wells were installed along White Oak Lane and south of the landfill in June 1999. Wells with screens at or near the water table were installed around the landfill in July 1999.

The HPA defines the zone of attenuation (ZA) around a landfill as being that area within 100 feet from the edge of the landfill and extending from the ground surface to the bottom of the uppermost aquifer, excluding the volume occupied by the waste. None of the monitoring wells at the facility are within this regulatory ZA. The downgradient wells range from 320 to 610 feet from the ZA, and the background wells are about 1670 feet from the ZA. It has been demonstrated that groundwater that has been affected by leachate has migrated beyond the regulatory ZA. Practically, the zone of attenuation is a swath between the waste and the Rock River. A map showing this existing ZA is attached. Well placement was determined based on the existing ZA, rather than the regulatory definition of ZA.

Well locations were specified based on the location of the potential source of contaminants (the CKD landfill), the assumed groundwater flow direction, and accessibility. The width of any potential contaminant plume downgradient from the landfill was established as the maximum width of the area containing CKD, normal to the groundwater flow direction. The groundwater flow direction was assumed to be from the landfill toward the Rock River, i.e., in a west-southwest direction. The maximum width of the landfill perpendicular to this direction, and the

width of any contaminant plume, was estimated to be 1550 feet wide. One set of wells, MW1-S and MW1-D, was located north-northeast of the landfill in the assumed upgradient location. These wells were intended to access groundwater that would be likely to travel beneath the landfill, but that had not yet been influenced by any potential leachate from the landfill. Wells MW2-S and MW2-D were located near the southern edge of the potential plume, and wells MW4-S and MW4-D were located near the northern edge of the potential plume. Wells MW3-S and MW3-D were located downgradient of the landfill and near the centerline of the potential plume, and they are in the existing zone of attenuation. Wells along White Oak Lane (MW5-S, MW5-D, MW6-S, MW6-D, MW7-S, MW7-D, MW8-S, and MW8-D) and between wells MW2 and MW3 were installed to better define the extent of groundwater affected by leachate. Wells screened at the water table were added to give a more complete three-dimensional picture of the groundwater quality.

Wells at each location are screened at different depths and formations. Wells designated with the suffix WT are screened at or near the water table, which is within the Pecatonica Formation near the landfill. Wells with the suffix S are screened in alluvium along White Oak Lane, or at the base of the Pecatonica elsewhere. Wells designated D are screened in the St. Peter Formation. One well with a suffix M was screened in the Glenwood Formation.

The St. Peter wells near the landfill were constructed by first drilling to the top of the first shale layer within the Glenwood Formation, and installing a 6-inch steel casing from the shale to the surface. The casing was grouted in place with high solids bentonite grout. This effectively scaled the base of the casing and prevented leakage of groundwater to lower aquifers. Drilling then proceeded inside the steel casing through the remaining Glenwood Formation and into the St. Peter Formation. Potable water was used for drilling fluid.

In all but one well, a 10 feot long PVC screen with 0.010-inch slots was used at the base of the borehole. The exception was well MW2-M, which has a 2 foot long screen. Centralizers were used on the screens in the deep wells to insure that the filter pack was of uniform thickness. Sand was placed around the screens up to a point 2 feet above the screens. About 3 feet of bentonite

pellets were then placed above the sand, and bentonite grout was tremied from the top of the pellets to the surface.

Well construction reports were completed for the monitoring wells on forms provided by the Illinois Department of Public Health (IDPH). These forms were submitted to IDPH, and copies are attached to this report.

The surveyed elevations of the wells and the static water levels are presented in the following tables.

## Table 3. Monitoring well and piezometer data. (all elevations are in feet above mean sea level)

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Table 4. Elevations of the potentiometric surface in monitoring wells.

(all elevations are in feet above mean sea level)

30-April-98	07-May-98	17-Jun-98	24-Jul-98	26-001-98	16-Feb-99
736.19	736.36	736.09	736.10	736.23	736.23
1.68.59	668.97	668.60	668.12	668.02	668.61
689.19	689.45	687.56	686.93	688.44	689.19
4-8-4.1-4	685.57	688 26	686.90	687.86	The state of the s
66246	662.44	10,130	(d(d) 10%	( (a) 4.4	(461.44
669.49	1009 20	68.08 24	KH (67.14	167.65	667.63
647.92	(547.60)	1945 Selector	645.59	\$546.01	646.23
683.87	683.63	680.59	677.35	682.10	682.33
657 85	657.70	1,5,5 () }	65324	655.30	655.83
ne de compression de la compression del compression de la compresi	n namade <mark>sana sana</mark> sanarang sebendar	703 80	703.36	703.37	703.60
	736.19 668.19 689.19 681.14 662.46 669.49 647.92	736.39 736.30 668.97 689.19 689.45 681.14 685.57 662.44 669.49 662.44 669.49 647.92 683.87 683.63	736.39 736.26 668.59 668.97 668.50 668.97 668.50 668.97 668.50 668.97 668.50 668.97 668.50 668.97 668.50 668.99 668.50 66	736.19         736.26         736.09         736.10           668.59         668.97         668.60         668.12           689.19         639.45         687.56         686.93           681.14         685.57         688.26         686.90           662.44         661.01         660.08           669.49         669.20         667.14           647.92         647.60         645.59           683.87         683.63         680.59         677.35           637.85         657.70         655.03         653.24	736.19         736.26         736.09         736.10         736.23           668.59         668.97         668.60         668.12         668.02           689.19         639.45         687.56         686.93         688.44           681.14         685.57         688.26         686.90         687.86           662.46         662.44         661.01         660.02         667.85           647.92         647.60         623.50         645.50         645.59         646.01           683.87         683.63         683.59         677.35         682.10           657.85         657.70         655.30         653.24         655.30

Propositionmentale sandines analysis for the St. Peter and Proposition Considerations are included in

#### 5.0 Croundwater Sampling

Prior to sompling, three well andmines of union were purged from each well with dedicated PVC besters. The grammal water semples were collected within 24 brown of purging the well. The amples are explained affirmally into the proper laboratory containers, and placed in a coorder with item. During the 1998 and 1999 example, amaples to be availyined for discolated metals were filtered in the field with simple-use 0.45 pm filters. A tandidity requirement was implemented during the flore appropriate and 2000. If the tembriday of the comple for medal analysis was greater than 5 NTO, the assumpter, one filtered and one filtered, were conferred. Samples that were filtered were intendified by the costain. "I'm other the most summe.

the 1998 and 1999, the unconficulty exclusive expension for activitie, every-arteside, and incorporate accordance and the exchangements. Made and the exchangements in our description in the description, and the exchangement, there are not described in all of the exclusion, along against every described in all of the exclusions are sent described in all of the exclusions are sentended in all of the exclusions are an entranced or supplies. Convertely there exclusions are exclusively there exists are exclusively there.

are thirty-four inorganic constituents, five volatile compounds, and pH that are being analyzed in the groundwater, leachate, and surface water.

#### 6.0 Groundwater Quality Standards

The groundwater quality standard for any constituent is the concentration of that constituent in the background groundwater. Wells MWI-WT, MWI-S, and MWI-D, screened at the water table in the uppermost aquifer (in the Pecatonica Formation), the base of the Pecatonica Formation, and the St. Peter Formation, respectively, are located upgradient from the landfill and are not expected to be infly—need by leachate from the landfill. Concentrations of constituents in the waters from these three wells represent background conditions, and were used to determine the groundwater quality standards.

Statistical tests were used on data from background water samples for the list of compounds that were analyzed at Dixon Marquette Ceznent Company. For most parameters there were four to fourteen sets of analytical data. The sampling frequency has not been great enough to establish if short term fluctuations in water quality occur, such as those that might occur because of rapid surface water influx, nor has it been long enough to show if variations recur seasonally. The population distribution was assumed to be normal.

Because of the small sample sizes, Student's I distribution was applied to sample sets where all the data were above the detection limit. In cases where 15 percent or less of the data was below the detection limit, the values below the detection limit were replaced with one-half of the detection limit. Student's I test was then applied. In cases where between 15 and 50 percent of the data was below the detection limit, Cohen's adjustment to the mean and variance was applied before Student's I test was applied. If 50 percent or more of the data was below the detection limit, a test of proportion will be used to determine if a downgradient well differs from the background well. For semi-volatile and volatile compounds, any detection is considered to be an increase above the background concentrations.

The concentrations of constituents that exceed the background concentrations at a level of significance of GOI were discounted. In cases where none of the constituent was detected in the

background well in any of the samples, any detection of the constituent in downgradient wells will be an exceedance of the background concentration. The following tables show each constituent and the minimum concentration that exceeds the background concentration, determined as described above, for the upper part of the uppermost aquifer (MW1-WT), the lower part of the uppermost aquifer (MW1-S), and the aquifer in the St. Peter Formation (MW1-D). Data in tables in Attachment 4 show which constituents in groundwater around the landfill exceed the background concentrations. Wells LW-1 and LW-3 were compared to well MW1-WT, and alluvial wells were compared to well MW1-S.

Table 5. Minimum Concentrations of Volatile Organic Compounds that Exceed Background Concentrations, in µg/l.

Parameter	kground Concent Water Table	Base of Pecatonica	Top of St. Peter
Acetone	50	50	50
Acrolein	50	50	50
Acrylonitrile	50	50	50
Benzene	5	5	5
Bromobenzene	5	5	5
Bromochloromethane	5	5	5
Bromodichloromethane	5	5	5
Bromoform	5	5	5
Bromomethane	10	10	10
n-butylbenzene	10	10	10
sec-butylbenzene	10	10	10
tert-butylbenzene	10	10	10
Carbon disulfide	10	10	10
Carbon tetrachloride	5	5	5
Chlorobenzene	5	5	5
Chlorodibromomethane	5	5	5
Chloroethane	10	10	10
Chloroform	5	5	5
Coloromethane	10	10	10
2-chlorotolnene	5	5	5
4-chlorotoluene	5	<u>\$</u>	5
Dibromomethane	\$	5	5
1,2-dichlorobenzene	5	5	5
1,3-dichlorobenzene	5	5	5
1,4-dichlorobenzene	5	.5	5
Dichlorodifluoromethane	10	1()	10
1,2-dichloroethane	5	5	5
1,1-dichloroethane	5	5	5
1,1-dichlorootherie	,5	5	5
cis-1,2 dichloroethene	5	5	5
trans-1,2-dichlemethene	5	5	5
trans-1.4-dichloro-2-butene	10	14)	1()
1,2-dichloropropasive	ŕ	5	\$
1,3-dichloropropane	3	5	5
2,2 dichloropropose	\$	5	<b>5</b> 1
1.1 dichloropropers	5	5	5
1,3-dichloropropene	10	10	10
cis-1,3-dichloropropene	5	Commission and the commission of the	R.
tomore I, 3-alochelessergeresgenese		Some or the contract of the co	5

Table 5. Minimum Concentrations of Volatile Organic Compounds that Exceed Background Concentrations, in μg/l.

Parameter	Water Table	Base of Pecatonica	Top of St. Peter
Ethylbenzene	5	5	5
2-hexanone	5	5	5
Hexachlorobutadiene	10	10	10
Iodomethane	10	10	10
Isopropylbenzene	10	10	10
p-isopropyltoluene	10	10	10
Methylene chloride	5	5	5
Methyl ethyl ketone	10	10	10
4-methyl-2-pentanone	5	5	5
Tetrahydrofuran	5.	5	5
Vinyl acetate	10	10	10
n-propylbenzene	10	10	10
Styrene	5	5	5
1,1,1,2-tetrachloroethane	5	5	5
1,1,2,2-tetrachloroethane	5	5	5
Tetrachloroethene	5	5	5
Toluene	5	5	5
1,2,3-trichlorobenzene	5	5	5
1,2,4-trichlorobenzene	5	\$	5
1,1,1-trichloroethane	5	5	5
1,1,2-trichloroethane	5	5	5
Trichloroethylene	5	5	5
Trichlorofluoromethane	10	10	10
1,2,3-trichloropropane	5	5	5
1,2,4-trimethylbenzene	10	10	10
1,3,5-trimethylbenzene	10	10	10
Vinyl chloride	10	10	10
o-xylene	10	10	10
m and p xylene	\$	5	5
Nylenes, total	5	5	5
Mote. Any Selection of a VOC' is a	considered to the americans	rice of the background concent	ration.

Table 6. Minimum Concentrations of Inorganic Compounds that Exceed Background Concentrations, in mg/l.

Parameter	Water Table	Base of Pecatonica	Top of St. Peter
Aluminum	0.454	0.04	(0.44)
Antimony	0.001	(0.11)	0.001
Arsenic	0.003	(0.10)	(0.18)
Barium	0.227	0.091	0.051
Beryllium	0.001	0.001	0.001
Boron	(0.43)	(0.30)	(0.20)
Cadmium	0.001	0.001	0.001
Calcium	126.1	109	91.8
Chromium	0.002	0.002	(0.36)
Cobalt	0.008	0.003	(0.30)
Copper	0.012	0 005	0.006
Iron	0.9	0.465	0.819
Lead	0.006	0.001	(0.18)
Magnesium	50.5	53.9	73.9
Manganese	2.69	0.045	0.098
Mercury	(0.14)	0.0002	0.0002
Nickel	0.023	0.007	0.003
Potassium	4.0	(0.50)	2.3
Selenium	(0.28)	0.004	0.005
Silver	0.001	0.001	0.001
Sodium	53.7	12.8	3.85
Thallium	(0.28)	(0.11)	(0.11)
Vanadium	0.007	(0.44)	0.003
Zinc	0.014	0.058	0.023
Alkalinity, total	342	253	346
Alkalinity, bicarbonate	334	250	343
Ammonia nitrogen	(0.14)	(0.20)	(0.22)
BOD, 5-day	54.5	(0.5)	(0.5)
Chloride	25.4	30.8	4.1
COD	(0.43)	5.0	5.0
Cyanide	0.1	0.005	0.005
Fluoride	0.76	(0.19)	0.19
Nitrate	(0.14)	11.07	0.21
Oil and Grease	(0.25)	(0.143)	0.1
Sulfate	242.8	178.3	44.2
Total dissolved solids	488	540	383
Toc	17	4.5	13.1

(Values in parentheses are the ratios of samples with detectable concentrations to the total number of samples, and are not background concentrations.)

Table 6. Minimum Concentrations of Inorganic Compounds that Exceed Background Concentrations, in mg/l.

Parameter	Water Table	Base of Pecatonica	Top of St. Peter
Aluminum	0.454	0.04	(0.44)
Antimony	0.001	(0.11)	0.001
Arsenic	0.003	(0.10)	(0.18)
Barium	0.227	0.091	0.051
Beryllium	0.001	0.001	0.001
Boron	(0.43)	(0.30)	(0.20)
Cadmium	0.001	0.001	0.001
Calcium	126.1	109	91.8
Chromium	0.002	0.002	(0.36)
Cobalt	0.008	0.003	(0.30)
Copper	0.012	0.005	0.006
Iron	0.9	0.465	0.819
Lead	0.006	0.001	(0.18)
Magnesium	50.5	53.9	73.9
Manganese	2.69	0.045	0.098
Mercury	(0.14)	0.0002	0.0002
Nickel	0.023	0.007	0.003
Potassium	4.0	(0.50)	2.3
Selenium	(0.28)	0.004	0.005
Silver	0.001	0.001	0.001
Sodium	53.7	12.8	3.85
Thallium	(0.28)	(0.11)	(0.11)
Vanadium	0.007	(0.44)	0.003
Zinc	0.014	0.058	0.023
Alkalinity, total	342	253	346
Alkalinity, bicarbonate	334	250	343
Ammonia nitrogen	(0.14)	(0.20)	(0.22)
BOD, 5-day	54.5	(0.5)	(0.5)
Chloride	25.4	30.8	4.1
COD	(0.43)	5.0	5.0
Cyanide	0.1	0.005	0.005
Fluoride	0.76	(0.19)	0.19
Nitrate	(0.14)	11.07	0.21
Oil and Grease	(0.25)	(0.143)	0.1
Sulfate	242.8	178.3	44.2
Total dissolved solids	488	540	383
TOC	17	4.5	5.1

(Values in parentheses are the ratios of samples with detectable concentrations to the total number of samples, and are not background concentrations.)

Table 7. Minimum Concentrations of Semivolatile Compounds that Exceed Back\_cound Concentrations, in ug/l. \*

Parameter	_cound Concent   Water Table	Base of Pecatonica	Top of St. Peter
Alachlor	1	1	1 .
Atrazine	<del>                                     </del>	<del></del>	<del>l i i</del>
Bis(2-ethylhexyl)phthalate	<del>                                     </del>	1	<u>                                     </u>
Bis(2-chloromethyl)ether		î	1
Di-n-butylphthalate	<del>1</del> 1	i	1
Diethyl phthalate	i	1	<del>l i</del>
Dimethyl phthalate	<del>                                     </del>	1	1
Isophorone	<del>                                     </del>	i	1
Naphthalene	1	1	1
p-cresol	1	i	1
Parathion	<del>                                     </del>	i	1
Pentachlorophenol	<del>1 1</del>	1	<u> </u>
Phenol		i	<del> </del>
Aldrin	0.034	0.034	0.034
gamma-BHC	0.025	0.025	0.025
Chlordane	0.037	0.037	0.037
4,4'-DDT	0.091	0.091	0.091
Dieldrin	0.044	0.044	0.044
Endrin	0.039	0.039	0.039
Heptachlor	0.040	0.040	0.040
Heptachlor Epoxide	0.032	0.032	0.032
Methoxychlor	0.176	0.176	0.176
Toxaphene	0.086	0.086	0.086
PCB-1016	0.050	0.050	0.050
PCB-1221	0.054	0.054	0.054
PCB-1232	0.065	0.065	0.065
PCB-1242	0.065	0.065	0.065
PCB-1248	0.090	0.090	0.090
PCB-1254	0.100	0.100	0.100
PCB-1260	0.100	0.100	0.100
PCB-1268	0.100	0.100	0.100
2,4-D	0.29	0.29	0.29
Silvex(2,4,5-TP)	0.34	0.34	0.34
Aldicarb	0.5	0.5	0.5
Carbofuran	5	5	5
1,2-dibromo-3-chloropropane	0.1	0.1	0.1
Ethylene dibromide	0.1	0.1	0.1

Note: A detection of a semivolatile compound in the downgradient monitoring wells is considered to be an exceedance of the background concentration.

### 7.0 Summary of Laboratory Analysis of Groundwater and Leachate

On May 7, 1998 the groundwater in the four shallow wells (MW 1-S through 4-S) and the four deep wells (MW 1-D through 4-D) was sampled. On June 16, July 24, and October 26, 1998, and February 1999, the groundwater in the same wells, and at a leachate well in the CKD landfill were sampled. In July, September, and November 1999 all wells except MW2-M were sampled, along with the spring near White Oak Lane. In 2000, sampling occurred in March, June, September, and November. The same locations were sampled in 2000 as in 1999.

The groundwater in the upgradient wells (MW1-S, MW1-D, and MW1-WT) had low concentrations of chloride, sulfate, potassium, and sodium. Dissolved solids were between 320 and 580 mg/l, which corresponds closely to values in Visocky et al. (1985) for the Galena-Platteville and Ancell Groups. The sample from the shallow aquifer had 8.2 µg/l toluene on July 24, 1998. The source of the toluene was not identified. In October 1998 and February 1999 no toluene was detected in MW1-S above 1 µg/l. Because toluene has not been confirmed in the well, its detection in the sample from July 1998 is considered to be sporadic, and toluene is not actually present in the well.

The leachate samples collected from LW1 and LW3 have elevated concentrations of calcium, potassium, sodium, chloride, sulfate, and total dissolved solids. In addition, the concentration of magnesium, manganese, iron are elevated when compared to the water table background concentrations of these parameters. Additional parameters that are slightly greater than the background concentrations include arsenic, boron, chromium, cobalt, copper, and lead. The leachate samples had low concentrations of benzene, 1,1-dichloroethane, and toluene.

The downgradient wells MW3-S and MW3-WT are located southwest of the landfill. The groundwater at this location has a pale yellow color and a high pH (>10). The analytical results show that the groundwater in both of these wells has elevated concentrations of chloride, sulfate, potassium, and dissolved solids when compared to the respective background concentration. In addition, elevated levels of sodium and alkalinity were seen in the shallow and water table wells at this location. Organic compounds were detected in these two wells (MW3-S and MW3-WT).

In MW3-5, three samples during 1998 had detected concentrations of 1,1-dichloroethane, the highest was 0.004 mg/L. The highest concentration of 1,1,1-trichloroethane in this well was 0.015 mg/L (15 µg/L) on May 7, 1998. In MW3-WT, 1,1-dichloroethane, 1,1-dichloroethane, and 1,1,1-trichloroethane were detected. The greatest concentration were 0.0081, 0.0079, and 0.010 mg/L, respectively.

Monitoring well 7-S is located along White Oak Lane, north of MW3. Groundwater at well MW7-S had a high pH, dissolved solids, potassium, sodium, chloride, and sulfate, which indicate contamination by leachate. For some sampling events the shallow wells along White Oak Lane exceeded the background concentrations of aluminum. arsenic, barium, boron, calcium, chromium, cobalt, copper, iron, lead, magnesium, manganese, nickel, potassium, sodium, vanadium, chloride, sulfate, dissolved solids, and alkalinity. The organic compounds 1,1-dichloroethane, 1,1-dichloroethene, and 1,1,1-trichloroethane were detected in MW7-S. The highest concentrations of the three compounds were 0.0639, 0.0082, and 0.006 mg/L, respectively.

The deep wells in this area (MW3-D and MW7-D) had elevated concentrations of barium, calcium, iron, manganese, nickel, potassium, sodium, alkalinity, and chloride exceeded the background concentrations. However, no organic compounds were detected during the sampling events.

Monitoring wells 5, 6 and 8 are also located along White Oak Lane, east of the landfill. In the shallow wells, elevated concentrations of aluminum, boron, calcium, copper, iron, manganese, potassium, sodium, alkalinity, chloride, sulfate and total dissolved solids were observed in these three locations. The groundwater in the St. Peter aquifer in this location is similar in composition.

The downgradient monitoring wells 2-S and 2-D, located south of the CKD landfill, had groundwater similar to the upgradient well. However, concentrations of barium, copper, magnesium, potassium, and alkalinity were slightly greater than the background concentration for these parameters in the shallow well.

Monitoring wells 9-S, 9-WT, and 9-D are located east of MW2 and south of the CKD landfill. In MW9-S and MW9-WT concentrations of aluminum, arsenic, calcium, magnesium, nickel, potassium, sodium, alkalinity, chloride, sulfate, and total dissolved solids consistently exceed the respective background concentrations.

Groundwater in the St. Peter aquifer at this location (MW2-D and MW9-D) has concentrations of barium, iron, manganese, potassium, chloride and sodium that are slightly greater than the background concentration in MW1-D. No organic compounds were detected in these monitoring wells.

The downgradient monitoring wells 4-S and 4-D, located northwest of the landfill, had groundwater with qualities intermediate between the upgradient well and the leachate well. The pH of the groundwater is 7.2. It had elevated concentrations of potassium, chloride, and sulfate, and somewhat elevated concentrations of sodium and dissolved solids. The current CKD waste disposal area is not directly upgradient from MW4-S, and there may be other waste disposal sites at the facility that have affected the groundwater at MW4-S. Previously there was a settling pond immediately north of MW4-S that received slurry and wastewater from a wet process in the #4 kiln. The #4 kiln was constructed in 1960. The exact dates of operation of the settling pond are not known, but a cover was being applied to it by 1977. Although this pond no longer exists, the disposal of process waste in it 1 ay have affected the groundwater.

Summary tables of the laboratory results have been included in attachment 4. The original laboratory reports are available from Preston Engineering, Inc. and will be provided on request.

### 8.0 Adjusted Groundwater Standards

Because the concentrations of several parameters in downgradient wells exceed the background concentrations and Section 620 class I groundwater standards, Dixon Marquette Cement Company is requesting adjusted groundwater standards. The following table lists the requested standards and the Class I groundwater standard.

Table 8. Requested Groundwater Standards all values in mg/L

	Class I Standard	Requested Adjusted Standard
Aluminum		2.522
Antimony	· 0.006	0.006
Arsenic	0.05	0.141
Barium	2	2
Beryllium	0.004	0.004
Boron	2	2
Cadmium	0.005	0.005
Calcium	***************************************	420
Chromium	0.1	0.1
Cobalt	1	j
Copper	0.65	0.65
Lead	0.0075	0.012
Iron	5	6.4
Magnesium		171.5
Manganese	0.15	3.3
Mercury	0.002	0.002
Nickel	0.1	0.1
Potassium		2309
Selenium	0.05	0.05
Sodium		140.5
Thallium	0.002	0.004
Vanadium		0.02
Zinc	5	5
Alkalinity, total		1699
Alkalinity, phenol.		1121
Alkalinity, bicarb		977
Ammonia nitrogen		1.25
Chloride	200	241
COD	•	187
Fluoride	4	4.4
Nitrate as N	10	11.1
pH (stnd units)		10.9
Sulfate	400	1245
TOC		29
Dissolved Solids	1200	4338
Benzene	0.005	0.005
1,1 dichloroethane		6.7
1,1-dichloroethene	0.007	35
Toluene	<u>1</u>	
1,1,1-trichlorocthane	0.2	0.2

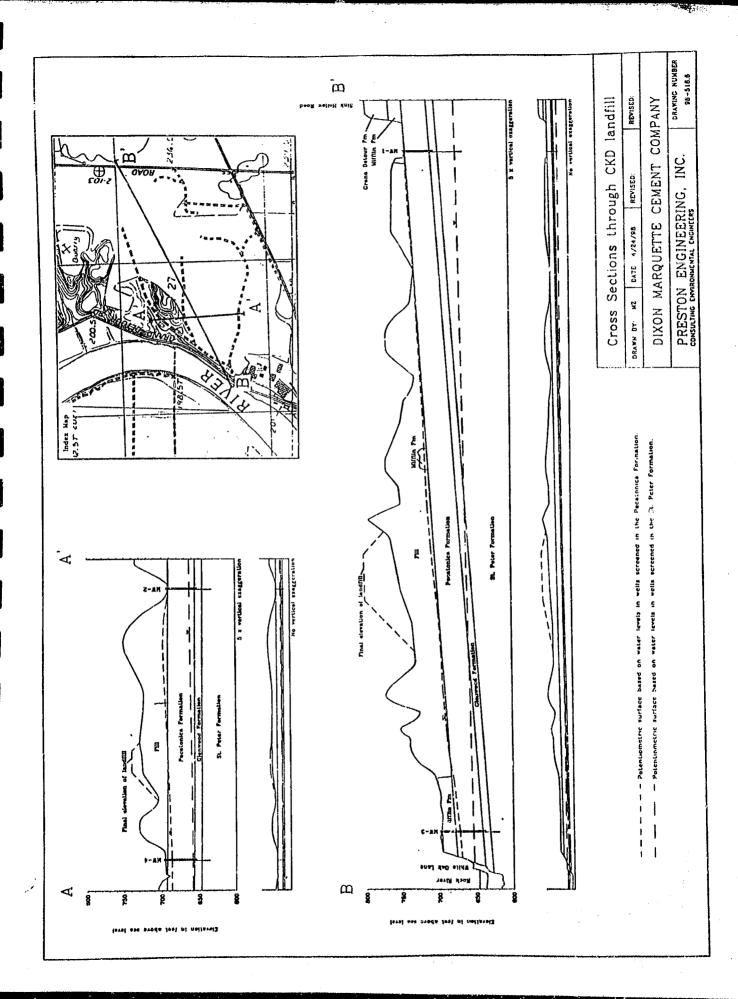
### 9.0 Abandonment of Borings and Monitoring Wells

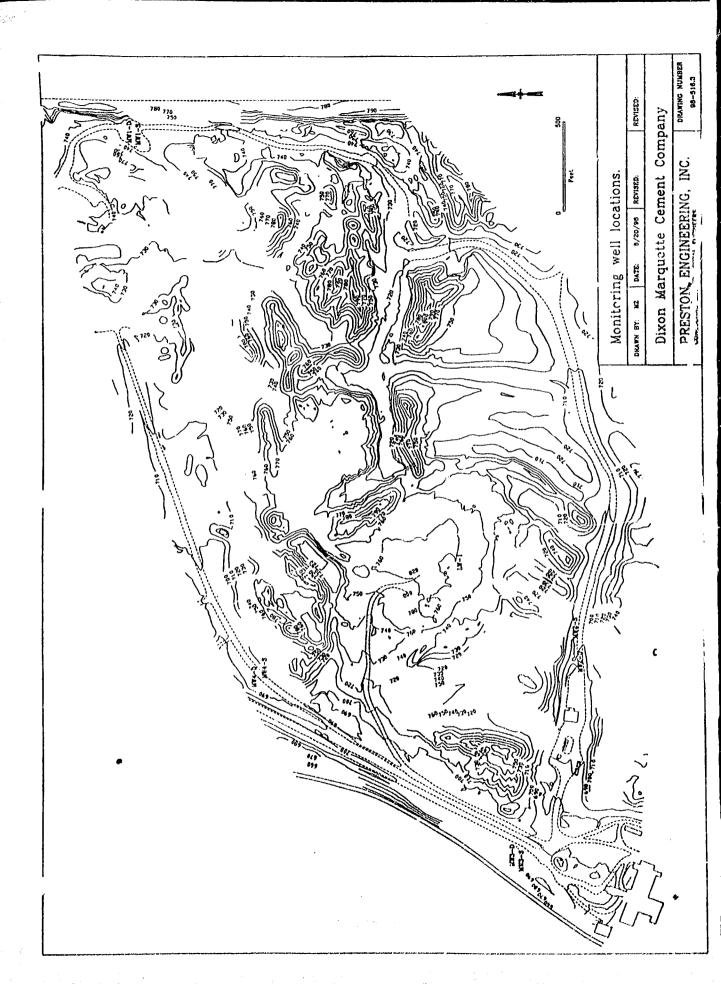
All borings made during the hydrogeologic investigation were converted to monitoring wells immediately following drilling. No borings have required abandonment.

Future borings that are not sealed or plugged immediately will be covered to prevent injury. Borings and wells that will be abandoned in the future will be filled with bentonite chips or pellets, or will have high-solids bentonite grout tremied into them. During tremie operations, the base of the tremie pipe will remain submerged in grout as the hole is being filled. Bentonite chips will be screened to remove fine material before being used. Bentonite and bentonite grout are compatible with the groundwater, leachate, soil, and CKD that are found at the site. Well casings that extend above the surface will be removed. All excess drilling mud, drill cuttings, and contaminated materials uncovered during or created by drilling will be disposed of in the on site landfill. Areas around the borings will be restored to conditions acceptable to Dixon Marquette Cement.

### Attachment 1

Cross Sections Through CKD Landfill





Attachment 2

Well Construction Reports

Date 5-5-98

FORM MUST BE COMPLETED WITHIN 30 DAYS OF COMPLETION AND SENT TO THE APPROPRIATE HEALTH DEPARTMENT TYPE OR PRESS FIRMLY WITH BLACK INK PEN, THIS

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- 2. Use: [] Domestic [] Irrigation [] Commercial [] Livestock Manitoring [] Other
- 3. Type of Well:
- qi, in. Depth 2. Borod Well: Hole Diameter
- in. Buried Slab: [ ]Yes [ ]No Casing Diameter
- b. Driven Well: Drive Pipe Diameter
  - in. Depth Casing Diameter 2 in Type PVC e. Drilled Well: Well Diameter\_

Cesing Grout:

Oversized

To(ft) Orill Hole(b) From(ft) Sentonite

S Z Gravel Pack: MYes Grain Size med, sand Finished In: Unconsolidated [ ] Rock 🔀

- 4. Well Disinfected? [ ]Yes [XNo
- none 5. Date Permanent Pump Installed
- 6. Licensed Pump Contractor\_ License Number
- 7. Pitless Adapter Installed? [ JYes PKNo Manufacturer

Attached to Casing - How? [] Screwed On [] Welded [] Compression

- 8. Type of Well Cap.
- gallons Captive Air: [ ]Yes [ ]No 9. Tank Working Cycle
  - 10. Pump and Equipment Disinfected? [ ]Yes [ ]No

General Comments: (If dry hole, fill out log & indicate how hole was sealed.)

Division of Environmental Health - 525 W. Jefferson Emois Department of Public Health

Springfield, IL 62761

DECOLARY ROTHER. The Same Agency is requesting disclosure of information that is necessary to accomplish the stendary purpose us confined under Public Act \$5.065. Disclosure of this information is mendatory. This orm has been secroted by the Forms Meascement Cotton.

GEULOGICAL AND WATER SURVEY WELL RECORD

- Date Issued 11. Permit Number
- Well # MW1-5 13. Drilling Capeny Name Geofechnical Services, Ixc. 12. Property Owner Dixon Marayette Cement Co.
  - 14. Name of Person who drilled the well Matt White
- Dixon, II 15. Well Site Address 1914 White Oak Lane.
- Land Dr 16. Twashp Name Dixon
- Elevation 739 Twashp 22 N Range 9 E 77 Sec 18. Location: Cuty Lee 17. Subdivision Name
  - Quarter Quarter of to. NE Quarter of the

20. Screen:

- Slot Size 0.010 Diameter 2 in. Material PVC Length 10 A. From(A) To (A) 228 19. Casing and Liner Pipe: Type 7 E 7
- ft. to 38 at depth 28 21. Water from Pecatorica Fm
- in. above ground level. bours. gpm for 22. Static Level 3.74 ft. below casing top which is 38 Pumping Level
- Depth Bottom(ft) Depth Top(ft) 이젊 23. Earth Materials Passed Through Pecatonica bolomite Sandshone Glenwood

Licensed Contractor Signature

Continue on back of sheet if necessary

License Number

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FORM MUST BE COMPLETED WITHIN 30 DAYS OF COMPLETION AND SENT TO THE APPROPRIATE HEALTH DEPARTMENT TYPE OR PRESS FIRMLY WITH BLACK INK PEN, THIS

- 1. Date Well Completed 3-20-48
- 2. Use: [] Domestic [] Irrigation [] Commercial [] Livestock [5] Monitoring [ ] Other
- 3. Type of Well:
- in. Buried Slab: [ ]Yes [ ]No in. Depth a. Bored Well: Hole Diameter
  - in. Depth b. Driven Well: Drive Pipe Diameter\_ Casing Diameter

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in. Depth c. Drilled Well: Well Diameter

Casing Dismeter 2 in Type

Joint

Casing Grout:	Oversized
Kind	Drill Hole(In) From(f)
Bentonite	525 66

. To(ft)

Gravel Pack: [XYes [ ]No Grain Size med, Sand Finished In: Unconsolidated [ ] Rock 🔀

Depth Bottom(ft)

Depth Top(R) gpm for

Pumping.

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Pumping Level

23. Earth Materials Passed Through

0

Lours.

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24 14.5

St. Peter sandstone Glenwood sandstone Pecatorica dolomite

- 4. Well Disinfected? [ JYes [XNo
- none 5. Date Permanent Pump Installed
- 6. Licensed Pump Contractor License Number
- 7. Pitless Adapter Installed? [ ]Yes [Mo

Attached to Casing - How? [ ] Screwed On [ ] Welded [ ] Compression Model Manufacturer

- 8. Type of Well Cap
- gallons Captive Air: [ ]Yes [ ]No 9. Tank Working Cycle
  - 10. Pump and Equipment Disinfected? [ ]Yes [ ]No

General Comments: (If dry hole, fill out log & indicate how hole was sealed.)

Division of Environmental Health - 525 W. Jefferson Illinois Department of Public Health Springfield, IL 62761

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CONTANT NOTICE. This State Agency is requesting disclosure of information that is necessary to a necessitish	sension years as outlined under Public Act \$5-0863. Disclosure of this information is mandatory.	anter.
5. This State Agency is reques	outlined under Public Act 15-	on has been approved by the Forms Managemont Contra
POICANT NOTICE	Managery purpose as	THE PART STORYED

Continue on back of sheet if necessary

Licensed Contractor Signature

### Geologic Descriptions

### Monitoring Well MW-1D (northeast side of CKD landfill)

(All measurements in feet)

Depth of Top	<u>Thickness</u>	<u>Unit</u>	<u>Description</u>
		Pecatonica Formation	
0	34		Limestone and dolomite
:			
34	44.5	Glenwood Formation	Sandstone and siltstone
••			
		St. Peter Formation	
44.5	78		Sandstone, quartz

## GEOLOGICAL AND WATER SURVEY WELL RECORD TYPE OR PRESS FIRMLY WITH BLACK INK PEN, THIS FORM MUST BE COLD!

2. Use: [] Domestic [] Irrigation [] Commercial [] Livestock [Monitoring [] Other 3. Type of Well:
[ ] Irrigation [ ] Commercial [ ] Live

		يه		
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in. Depth	in. Buried Slab: []Yes []No	ia. Depth	in. Depth	Joint
		b. Driven Well: Drive Pipe Diameter	c. Drilled Well: Well Diameter 2	r in Type
2. Bored Well: Hole Diameter	Casing Diameter	b. Driven Well:	c. Drilled Well:	Casing Diameter

Gravel Pack: NYes []No Grain Size ned Finished In: Urconsolidated [] Rock N

- 4. Well Disinfected? [ ]Yes [MNo
- 5. Date Permanent Pump Installed
- 7. Pitless Adapter Installed? [ ]Yes MNo 6. Licensed Pump Contractor\_ License Number
- Attached to Casing How? [ ] Screwed On [ ] Welded [ ] Compression Model Manufacture

8. Type of Weil Cap.

gallons Captive Air: [ ]Yes [ ]No 10. Pump and Equipment Disinfected? [ ]Yes [ ]No 9. Tank Working Cycle

General Comments: (If dry hole, fill out log & indicate how hole was sealed.)

Continue on back of sheet if necessary

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MOSTANT NOTICE. This Same Agency is requesting disclosure of information that is necessary to accompli	se sunderly purpose as outlined under Public Act 15-0663. Disclosure of this information is mendatory. This	arm has been morewed by the Forms Manacoment Center.	
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MORTANT NOTICE	to sestions purpose as on	arm has bren anoroved to	

20. Well # MW2-5 es Inc. Xan IL Land ID#	Elevation 692 flows. N. Range 9 E. Our et 20. Screen:	Diameter Zin. Length 10 ft. Slot Size 0.010 Material PVC	ch is 28 in above ground level.  gpm for bours.  epth Top(ft) Depth Bottom(ft)	34.5	
Date Issued  E Coment Co.  ical Services  natt White  ak Lane. Dixa	Sect 27 Twishp 22 N Range 9 E SE Quarter of the 51 V Qur. or 20. Screen:	24,5 O	at depth 34.5 which is 28 gpm for Depth Top(ft)	36.5	
11. Permit Number  12. Property Owner Dixon Manyvette Cennent Co. Well # MW2-S  13. Drilling Company Name Sected high Services Inc.  14. Name of Person who drilled the well Matt White  15. Well Site Address 1914 White Oak Lane Dixon IL  16. Twisto Name Dixon  Land ID#	1 1 1 1	·	21. Water from Pecarlonica Fran at depth 3. 22. Static Level 3.8 ft. below casing top which is 2.8 Pumping Level ft. Pumping gpm for 23. Earth Materials Passed Through Depth Top(ft)	recatonica polomite Glenwood Sandstane	

Licensed Contractor Signature

## WELL CONSTRUCTION REPORT

Date 5-5-48

W-

FORM MUST BE COMPLETED WITHIN 30 DAYS OF COMPLETION AND SENT TO THE APPROPRIATE HEALTH DEPARTMENT (13) 7-TYPE OR PRESS FIRMLY WITH BLACK INK PEN, THIS

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Commercial	
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Irrigation	] Other
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. Use: [ ] Domestic	[N. Monitoring
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Depth	[ ]Yes [ ]No
ij	Buried Slab:
Hole Diameter	ter in.
a. Bored Well:	Casing Diame

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	Diameter
	Pipe 1
	Drive
	Well:
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Depth 43,54. .넘 ii. Type PVC c. Drilled Well: Well Diameter

Diameter 4	
Cesing 1	Casing Grout:

9 (1011)	O CLISICAL CONTRACTOR		
Kind	Drill Hole(In) From(ft), To(ft)	From(ft)	, To(ft)
Bentonite	4.25	0 h	0

Gravel Pack: [MYes [ ]No Grain Size med, SANO Finished In: Unconsolidated [ ] Rock 🔀

- 4. Well Disinfected? [ ]Yes DANo
- 5. Date Permanent Pump Installed

6. Licensed Pump Contractor License Number

7. Pitless Adapter Installed? [ ]Yes [MNo Manufacturer

Model

Attached to Casing - How? [] Screwed On [] Welded [] Compression

- 8. Type of Well Cap
- Captive Air: [ ]Yes [ ]No gallons 9. Tank Working Cycle
- 10. Pump and Equipment Disinfected? [ ]Yes [ ]No

General Comments: (If dry hole, fill out log & indicate how hole was sealed.)

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Division of Environmental Health - 525 W. Jefferson Springfield, IL 62761 DATORITANT NOTICE. This Same Agency is requesting disclosure of information that is necessary to accomplish the sensiony purpose as outlined under Public Act \$5.0863. Disclosure of this information is mendatory. This stoved by the Forms Management Contex. form has been me

GEOLOGICAL AND WATER SURVEY WELL RECORD	ORD
11. Permit Number Date Issued	sued
12. Property Owner Dixon Marginette Cement Co. Well # MW2-	Well # MW2-
13. Drilling Company Name Geofechnical Services Inc	ces Inc
14. Name of Person who drilled the well Matt White.	
15. Well Site Address 1914 White Oak Lane, Dixon, IL	Dixon, IL
16. Twastp Name Dixon	Land ID#
17. Subdivision Name Lot	Elevation
ee Sect 27 Tw	22 N Range 9E
SE	V Quarter
19. Casing and Liner Pipe:	20. Screen:
Dia (In) Type From(ft) To (ft)	Diameter 2 in.
PVC	Length 2 ft.
	Slot Size 0.010
	Material PVC
21. Water from Glenwood Fro at depth 4	at depth 41,5 A. to 43,5 A.
22. Static Level 8,71 ft. below casing top which is 28	in. above ground level.
Pumping Level ft. Pumping gpm for	hours.
assed Through	Depth Top(ft) Depth Bottom(ft)
	36.5
Glenwood sandstone 36.5	43,5

Continue on back of sheet if necessary

Licensed Contractor Signature

License Number

GEULOGICAL AND WATER SURVEY WELL RECORD

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- 3-20-48 1. Date Well Completed
- 2. Use: [] Domestic [] Irrigation [] Commercial [] Livestock [X Monitoring [ ] Other\_
- 3. Type of Well:
- in. Depth a. Bored Well: Hole Diameter
- in. Buried Slab: [ ]Yes [ ]No Casing Diameter
- <u>ن</u>ه 90 in. Depth b. Driven Well: Drive Pipe Diameter
  - in. Depth\_ 2 in Type c. Drilled Well: Well Diameter Casing Diameter

Casing Grout:

Oversized

To(ft) Drill Hole(In) From(ft) % ₩ 5.75 Bentonite

Gravel Pack: [NYes []No Grain Size med, Sand Finished In: Unconsolidated [ ] Rock [3

- 4. Well Disinfected? [ ]Yes ONNo
- 5. Date Permanent Pump Installed
- 6. Licensed Pump Contractor
  - License Number
- Attached to Casing How? [ ] Screwed On [ ] Welded [ ] Compression 7. Pitless Adapter Installed? [ ]Yes MNo Manufacturer
  - 8. Type of Well Cap.
- gallons Captive Air: [ ]Yes [ ]No 9. Tank Working Cycle
  - 10. Pump and Equipment Disinfected? [ ]Yes [ ]No

General Comments: (If dry hole, fill out log & indicate how hole was scaled.)

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Division of Environmental Health - 525 W. Jefferson Springfield, IL 62761 DEFORMANT NOTICE. This State Agency is requesting disclosure of information that is necessary to accomplish the strategy propose as cottless under Public Act 35-0863. Disclosure of this information is mendelony. This form has been sessived by the Forms Menament Center.

11. Permit Number	Date Issued	
12. Property Owner Dixon Margy ette Cement Co Well # MINZ - D	e Cement Co	Well # MN2-D
13. Drilling Company Name Geofechnical Services. Inc.	cal Services.	Lnc.
14. Name of Person who drilled the well Matt White.	Matt White	
15. Well Site Address 1914 White Oak Lane Dixon	at Lane Dix	n II
16. Twishp Name Dixary	נ	₹
17. Subdivision Name	ទ	Elevation 692, ft.
18. Location: Cnty Lee Sect 27		Twash 22 N Range 9E
of the SE		Quarter
		20. Screen:
Dia (In) Type	From(ft) To (ft)	Diameter 2 in.
o u	50 0	Length 10 ft.
		Slot Size 0,010
		Material PVC
21. Water from St. Poller Fr	at depth 2	ft. to 60 ft.
22. Static Level 30,45ft. below easing top which is 27		in. above ground level.
Pumping Level ft. Pumping	gpm for	hours.
Passed Thro	Depth Top(ft)	Depth Bottom(ft)
	0	R N
Gleninood sendstone	36.5	エ
St. Peter Sandstone	4	3

Licensed Contractor Signature

Continue on back of sheet if necessary

License Number

### Geologic Descriptions Monitoring Well MW-2D (south side of CKD landfill) (All measurements in feet)

Depth of Top	Thickness	Unit	Description
0	1.5		Fill material, crushed rock
·	. 1	Pecatonica Formation	
		Medusa member	
1.5	4.5		Limestone, yellow with gray dolomitic mottling, rare brachiopod
		New Glarus member	
6	1		Limestone, yellow with dark yellow mottling
7	4.5		Limestone, yellow with gray dolomitic mottling, some brachiopods, gastropods, crinoids
11.5	4.5		Limestone, yellow and gray, stylolites at 12 feet
		Dane member	
16	7.5		Limestone, yellow with large dolomitic or shaly mottling, some brachiopods, pyritic at 19.5 feet, calcite in vug at 22 feet
23.5	4		Dolomite, gray, sandy, vuggy from 25 to 26 feet and 27 to 27.5 feet, calcite in vug at 24 feet, some pyrite
27.5	1.5		Limestone, yellow and gray with dolomitic mottling
		Chana member	·
29	7.5		Dolomite, gray, sandy, pyritic, vuggy at 32 to 35.5 feet
		Glenwood Formation	
36.5	2.5		Sandstone, fine grained quartz, some pyrite
39	0.5		Siltstone, sandy, gray, some pyrite
39. <b>5</b>	1.5		Siltstone, gray, fissile
41	2		Sandstone, quartz, some pyrite
43	1		Siltstone, gray, fissile
44	3		Sandstone, quartz, mottled, some pyrite
		St. Peter Formation	
47	2.5		Sandstone, quartz, poorly cemented, pyritic at 47 feet
49.5	10		Sandstone, quartz

# TYPE OR I FORM MU

TE OR PRESS FIRMLY WITH BLACK INK PEN, THIS	NRM MUST BE COMPLETED WITHIN 30 DAYS OF COMPLETION	ID SENT TO THE APPROPRIATE HFALTH DEPARTMENT	Date Well Completed 3-20-48	Use: [] Domestic [] Irrigation [] Commercial [] Livestock	Da Monitoring [ ] Other
TH OK PKE	RM MUST E	D SENT TO	Date Well C	Use: [] Don	IX Mon

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	in. Depth	in. Buried Slab: [ ]Yes [ ]No	ter in. Depth	Z in Derth	Joint	Oversized
	a. Bored Well: Hole Diameter		b. Driven Well: Drive Pipe Diameter_	c. Drilled Well: Well Diameter Z in. Depth	ter in Type	
3. Type of Well:	a. Bored Well:	Casing Diameter	b. Driven Well:	c. Drilled Well:	Casing Diameter	Casing Grout:

Kind	Drill Hole(In) From(ft), To(ft)	From(ft)	. To(ft)
Bentonite	9)	4	4
Contact of Transport Stated ( 1 Contact Dark: DXVec   1No	Speniel Dack. CXV	Ne l'INo	

Gravel Pack: [Ayes | JNo Grain Size Med, Sand Rock [X

S N
[]Yes
Disinfected?
Well
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Installed
Pump
Permanent
Date
'n

		[]Yes DNo
6. Licensed Pump Contractor	Vumber	7. Pitless Adapter Installed? [ ]Yes Offlo
6. Licensed	License Number	7. Pitless A

	] Compression
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gallons Captive Air: [ ] Yes [ JNo	
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9. Tenk Working Cycle	10. Pump and Equipment
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General Comments: (If dry hole, fill out log & indicate how hole was sealed.)

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linois Department of Public Health	Nision of Environmental Health - 525 W. Jefferson	
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CE. This State Agracy is requesting di	as outlined under Public Act 15-0863.	form has been approved by the Forms Manarement Conter.
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<b>a</b>	well # MW3-S Inc	CON IL	Elevation	Twish 22 N Range 4E fithe SW Quarter	20. Screen:	Diameter 7 in.	Length 10 ft.	Material PVC	ft. to 46 ft.	in. above ground level.	hours.	Depth Bottom(ft)	1.5	18.5 18.5	55	g.					
WELL RECOR	Ement Co.	K Lane Dixo	Lot			From(ft) To (ft)	8		at depth 59		gpm for		0	1.5	18.5	55					
GEOLOGICAL AND WATER SURVEY WELL RECORD	11. Permit Number  12. Property Owner Dixon Macayette Cement Co. Well is Drilling Company Name Geofechylcal Services Inc. 13. Name of Person who drilled the well Matt White	15. Well Site Address 1914 White Oak Lane Dixon IL	17. Subdivision Name	of the NW	ļ.				21. Water from Pecatonica Fm	22. Static Level 26.80ft. below casing top which is 28	Pumping Level ft. Pumping	Passed	Ş	Mifflin dolomite	Pecatanica dolomite	Gleningod Sandstone					

Continue on back of sheet if necessary

Licensed Contractor Signature

Date 55-98

Date Issued

GEOLOGICAL AND WATER SURVEY WELL RECORD

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AND SENT TO THE APPROPRIATE HEALTH DEPARTMENT	11. Permit Number

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- 3. Type of Well:
- in. Depth a. Bored-Well: Hole Diameter
- in. Buried Slab: [ ]Yes [ ]No Casing Diameter
- in. Depth b. Driven Well: Drive Pipe Diameter
  - in. Depth 78 Casing Diameter 2 in. Type FVC 11 c. Drilled Well: Well Diameter

Casing Grout:

ng digni:	00000000		
Kmd	Drill Hole(In) From(ft), To	From(ft)	7
Bentonite	5.25	99	13

%[] Gravel Pack: [AYes Finished In: Unconsolidated [ ]

Grain Size Med, Sand Rock 💢

- 4. Well Disinfected? [ ]Yes DANo
- 5. Date Permanent Pump Installed
- 6. Licensed Pump Contractor\_ License Number
  - 7. Pitless Adapter Installed? [ ]Yes MNo

Attached to Casing - How? [ ] Screwed On [ ] Welded [ ] Compression Manufacturer

- 8. Type of Well Cap
- gallons Captive Air: [ ]Yes [ ]No 9. Tank Working Cycle
  - 10. Pump and Equipment Disinfected? [ ]Yes [ ]No

General Comments: (If dry hole, fill out log & indicate how hole was sealed.)

Health	
Public	
nt of	
Departme	,
linois	

Division of Environmental Health - 525 W. Jefferson

Springfield, IL 62761

ILTORITANT NOTICE. This Suce Areacy is requesting disclosure of information that is necessary to accomplish the state of purpose as ordined under Public Act \$5-0663. Disclosure of this information is mandatory. This form has been arouved by the Forms Manatement Center.

Cement Co Well # MW3-D	at White	ak Lane. Dixon II.	Land ID#	Lot Elevation ft.	Sect 27 Twishp 22 N Range 9E	NW Quarter of the SW Quarter	20. Screen:	From(ft) To (ft), Diameter Z in.	68 0 Length 10 ft.	Slot Size 0.0 0	Material PVC	at depth 68 ft. to 78 ft.		gpm for hours.	Depth	0 (,5	1.5 18.5	18.5 55	55 66	66 78					
12. Property Owner Dava Manyette Cement Co Well # MW3-D	13. Drilling Company Name Geotechnical Services Lnc	15. Well Site Address 1914 White On	16. Twarip Name Dixon	17. Subdivision Name	ee	MΝ	19. Casing and Liner Pipe:		Lin PVC			21. Water from St. Peter Fr	_	Pumping Level ft. Pumping	Passed Thro	(8)	Mifflin dolomite	Pecatonica dolomite	Glenwood sandstone	St. Peter sandstone					

Continue on back of sheet if necessary

Licensed Contractor Signature

11/95

### **Geologic Descriptions**

### Monitoring Well MW-3D (southwest side of CKD landfill)

(All measurements in feet)

Depth of Top	Thickness	Unit	Description
0	1.5		Fill material, crushed rock
		*.	
		Mifflin Formation	
1.5	18.5		Limestone
			**
		Pecatonica Formation	
18.5	55		Limestone and dolomite
		Glenwood Formation	
55	66		Sandstone and siltstone
		St. Peter Formation	
66	78		Sandstone, quartz

Date

Well # MW 4-5

YPE OR PRESS FIRMLY WITH BLACK INK PEN, THIS	GEOLOGICAL AND WATER SURVEY WELL RECORD
ORM MUST BE COMPLETED WITHIN 30 DAYS OF COMPLETION	
ND. SENT TO THE APPROPRIME HEALTH DEPARTMENT CONTRACT TO THE APPROPRIME Number	11. Permit Number

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- 2. Use: [ ] Domestic [ ] Irrigation [ ] Commercial [ ] Livestock Monitoring [] Other
- 3. Type of Well:
- in. Buried Slab: [ ]Yes [ ]No in. Depth 4. Bored Well: Hole Diameter\_ Casing Diameter
- in. Depth 5. Driven Well: Drive Pipe Diameter\_
- c. Drilled Well: Well Diameter 4 20 Zm. Deptt 42. ft. Joint is. Twe PVC Casing Diameter 2

sized	Drill Hole(In) From(ft)	25 30		
Oversized	Drill	52.h		
Casing Grout:	Kind	Bantonife		

To(ft)

Gravel Pack: . [MYes [ ]No Grain Size med, Sand Finished In: Unconsolidated [ ] Rock 🔀

- 4. Well Disinfected? [ ]Yes [MNo
- 5. Date Permanent Pump Installed Aone
- 6. Licensed Pump Contractor License Number
- Attached to Casing How? [ ] Screwed On [ ] Welded [ ] Compression Model 7. Pitless Adapter Installed? [ ]Yes [Mo Manufacturer
- 8. Type of Well Cap
- gallons Captive Air: [ ]Yes [ ]No 10. Pump and Equipment Disinfected? [ ]Yes [ ]No 9. Tank Working Cycle

General Comments: (If dry hole, fill out log & indicate how hole was sealed.)

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linois Department of Public Health hivision of Environmental Health - pringfield, IL 62761	
linois Department of hivision of Environme pringfield, IL 62761	
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g discionary of information that is necessary to accomplish the sametory purpose as sufficiel pusher Public Act 15-0863. Discionary of this information is m form has been experend by the Forms Menarement Creter IMPORTANT NOTICE. The Same Again, is me

N			- {	ë g									•												-
Well # MW 4-		Dixon, II-	Land ID#	Elevation (86	Twash 22 N Range 9 E	/ Quarter	Ħ	Diameter 2 in.	Length 10 ft.	Slot Size 0.010	Material PVC	A. to 42 A.	in. above ground level.	bours.	Depth Bottom(ft)	8.5	38.5	777							
Cement Co.	Matt White	ak Lane. D		Ĭ	27 Twishp 22			From(ft) To (ft)	32 0			at depth 32	p which is 28	gpm for	Depth Top(ft)	0	8.5	38.5							
12. Property Owner Dy you Macquette Cement Co. Well # MW 4-5	14. Name of Person who drilled the well Matt White	15. Well Site Address 1914 White Oak Lane,	16. Twishp Name Dixon	17. Subdivision Name	Sect Sect	SE	19. Casing and Liner Pipe:	Dia (In.) Type	á		•	21. Water from Pacatonica Em	22. Static Level 1090ft. below casing top which is 28	Pumping Level ft. Pumping	Passed Thro	Soil	Pecatonica dolomita	Glenwood sandstone							Continue on back of sheet if necessary

FORM MUST BE COMPLETED WITHIN 30 DAYS OF COMPLETION. AND SENT TO THE APPROPRIATE HEALTH DEPARTMENT TYPE OR PRESS FIRMLY WITH BLACK INK PEN, THIS

2000	21-07-4
	1 Completed
	Date Well

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[ ] Irrigation	7 C L
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a. Bored Well: Hole Diameter_	ब्र	b. Driven Well: Drive Pipe Diameter
Well:	Casing Diameter	Well:
Bored	Casing	Drives
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Depth 6	Joint
Well Diameter Z in.	7 in Type PVC
c. Drilled Well: V	Cacino Diameter

n. Type PVC	Oversized
Casing Diameter 2 in. T	Casing Grout

	(ft) To(ft)	7 5.84		
Oversized	Drill Hole(In) From(ft), To(ft)	5.25 48		
ر				
ing Grout	Kind	Semborte		

Gravel Pack: NYes [ INo Grain Size Mech Sand Finished In: Unconsolidated [ ]

4. Well Disinfected? [ ]Yes DNo

Rock [X

none 5. Date Permanent Pump Installed

6. Licensed Pump Contractor\_ License Number

Attached to Casing - How? [ ] Screwed On [ ] Welded [ ] Compression 7. Pitless Adapter Installed? []Yes PANo Manufacturer

8. Type of Well Cap.

gallons Captive Air: [ ]Yes [ ]No 9. Tank Working Cycle

10. Pump and Equipment Disinfected? [ ]Yes [ ]No

General Comments: (If dry hole, fill out log & indicate how hole was scaled.)

Illinois Department of Public Health

Division of Environmental Health - 525 W. Jefferson Springfield, IL 62761 IMPORTANT NOTICE. This Same Agency is requesting disclosure of information that is mecessary to accomplish the standard purpose as outlined under Public Act \$5.065. Disclosure of this information is mendatory. This form has been scorered by the Forms Maintenant Center.

## GLULOGICAL AND WATER SURVEY WELL RECORD

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Date Issued	, ,
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Matt White	Lane Dixon, IL
14. Name of Person who drilled the well Mat-	15 Well Site Address 1914 White Oak Lane Dixon.
14. Name of Person w	15 Well Site Address

77 N Range 9 E	1
me Lot Elevation 60	I Cheman
17. Subdivision Name Lot	, o Table 10 10 10 10 10 10 10 10 10 10 10 10 10

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19 Casing	19 Casing and Liner Pipe:		20. Screen:
رق) وزر	Type	From(ft) To (ft),	Diameter (- III.
7	7 is PVC		Length 10 ft.
			Slot Size 0.0/C
			Material PVC
		1.1. KO.F.	1. 1. 1. 1. 1. 1. 1. 1. 1. 1. 1. 1. 1. 1

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22. Static Level 37,2(At. below casing top which is 29 in. above	Towns I out

							 	 	 	 	-
pours.	Depth Bottom(ft)	8,5	38.5	48,5	60.5						
gpm for	Depth Top(ft)		8,5	38.5	5'8h						
Pumping Level ft. Pumping	send Through		Permiss dolowite	Cleawood Sand Stane	St. Dator candstone						

Continue on back of sheet if necessary

Licensed Contractor Signature

License Number

11/95

Attachment 3

Aquifer Test Data



June 12, 1998

Mr. Mark Zell
Preston Engineering
4436 N. Brady Street
Davenport, Iowa 52806-4009

RE: PACKER TEST RESULTS

**DIXON MARQUETTE CEMENT COMPANY** 

DIXON, ILLINOIS

Dear Mr. Zell:

Geotechnical Services, Inc. (GSI) has completed the analysis of the packer tests conducted at the above referenced facility. The results are presented in the attachments to this letter. The following equation was used for calculating the hydraulic conductivity of the test intervals.

$$K := \frac{O}{2 \cdot \pi \cdot L \cdot H} \cdot \operatorname{asinh} \left( \frac{L}{2 \cdot r} \right)$$

(Based on equation from Earth Manual, US Dept. of Interior, Bureau of Reclamation)

K = hydraulic conductivity

Q = constant rate of flow into the portion of borehole

L = length of portion of the borehole tested

H<sub>g</sub> = head due to gravity

 $H_p$  = head as applied pressure

H = differential head of water

r = radius of hole tested

sinh<sup>-1</sup> = inverse hyperbolic sine

Please note that on three of the tests, MW-1, MW-3 and MW-4) no flow was measured during the packer tests. To provide a maximum hydraulic conductivity estimate, the flow meter minimum detection of 0.0017 gpm was used as the flow rate for these tests. This allows for a maximum hydraulic conductivity estimate to be calculated.

GEOTECHNICAL, MATERIALS ENGINEERING & ENVIRONMENTAL CONSULTANTS

258 EAST 90TH STREET DAVENPORT, KOWA 52808-7341 (319) 285-8541 • FAX (319) 285-8545



If you have any questions please call.

Sincerely, GEOTECHNICAL SERVICES, INC.

Joel P. Zirkle, ICGP, RG Senior Geologist

cc: File 26Q0026

Attachments



### Attachment A

Date Test Conducted: March 31, 1998

Location: MW-1

Depth of Test Interval: 42.5 ft. to 43.5 ft.

$$r := \frac{1.5 \text{ in}}{12 \cdot \frac{\text{in}}{\text{ft}}}$$

$$Q = 0.0017 \frac{\text{gal}}{\text{min}} \cdot 0.13368 \frac{\text{ft}^3}{\text{gal}}$$

Hp := 35 psi 2.31 
$$\frac{\text{ft}}{\text{psi}}$$

$$H = Hg + Hp$$

$$L = 1 \cdot ft$$

$$K = \frac{Q}{2 \cdot \pi \cdot L H} \cdot \operatorname{asinh} \left( \frac{L}{2 \cdot r} \right)$$

$$K = 5.83 \cdot 10^{-7} \cdot \frac{ft}{min}$$

$$K = 2.9 \cdot 10^{-7} \cdot \frac{\text{cm}}{\text{sec}}$$



### Attachment B

Date Test Conducted: March 24, 1998

Location: MW-2

Depth of Test Interval: 42.0 ft. 10 43.5 ft.

$$r := \frac{1.5 \text{ in}}{12 \cdot \frac{\text{in}}{\text{ft}}}$$

$$Q = 0.230 \frac{\text{gal}}{\text{min}} \cdot 0.13368 \frac{\text{ft}^3}{\text{gal}}$$

$$Hg = 48.5 ft$$

Hp := 20 psi 2.31 
$$\frac{ft}{psi}$$

$$H = Hg + Hp$$

$$K = \frac{Q}{2 \cdot \pi \cdot L H} \cdot a \sinh \left( \frac{L}{2 \cdot r} \right)$$

$$K = 8.58 \cdot 10^{-5} \cdot \frac{ft}{min}$$

$$K = 4.36 \cdot 10^{-5} \cdot \frac{\text{cm}}{\text{sec}}$$



### Attachment C

Date Test Conducted: March 27, 1998

Location: MW-3

Depth of Test Interval: 62.0 ft. to 63.0 ft.

$$r := \frac{1.5 \text{ in}}{12 \cdot \frac{\text{in}}{\text{ft}}}$$

Q = 0.0017 
$$\frac{\text{gal}}{\text{min}} \cdot 0.13368 \frac{\text{ft}^3}{\text{gal}}$$

$$Hg := 68.5 \text{ ft}$$

Hp := 50 psi 
$$\cdot$$
 2.31  $\cdot \frac{\text{ft}}{\text{psi}}$ 

$$H := Hg + Hp$$

$$L = 1.0 \text{ ft}$$

$$K = \frac{Q}{2 \pi \cdot L H} \cdot a \sinh \left( \frac{L}{2 r} \right)$$

$$K = 4.12 \cdot 10^{-7} \cdot \frac{ft}{min}$$

$$K = 2.09 \cdot 10^{-7} \cdot \frac{cm}{sec}$$



### Attachment D

Date Test Conducted: April 3, 1998

Location: MW-4

Depth of Test Interval: 45.0 ft. to 46.0 ft.

$$r := \frac{1.5 \text{ in}}{12 \cdot \frac{\text{in}}{\text{ft}}}$$

$$Q := 0.0017 \frac{\text{gal}}{\text{min}} \cdot 0.13368 \frac{\text{ft}^3}{\text{gal}}$$

$$Hg := 51.5 ft$$

$$Hp := 40 \text{ psi} \cdot 2.31 \cdot \frac{\text{ft}}{\text{psi}}$$

$$H := Hg + Hp$$

$$L := 1 \cdot ft$$

$$K = \frac{Q}{2 \cdot \pi \cdot L \cdot H} \cdot \operatorname{asinh} \left( \frac{L}{2 \cdot r} \right)$$

$$K = 5.27 \cdot 10^{-7} \cdot \frac{ft}{min}$$

$$K = 2.67 \cdot 10^{-7} \cdot \frac{cm}{sec}$$

### Slug Test Results

Slug out Title: Client: DMC 98-516 Job Number:

Well Number: MW-1 Shallow

### Well Geometry

38.45 H: 12. Le: 38.45 Lw: .083 rc: 3. rw:

0.0 filter pack porosity: effective radius: 8.33E-2

### Bouwer Rice Coefficients

Le/rw: 4. 1.705 A: .201 B: C: .71 ln(Re/rw): 1.643

Hvorslev

52.228 F:

Least Squares Fit

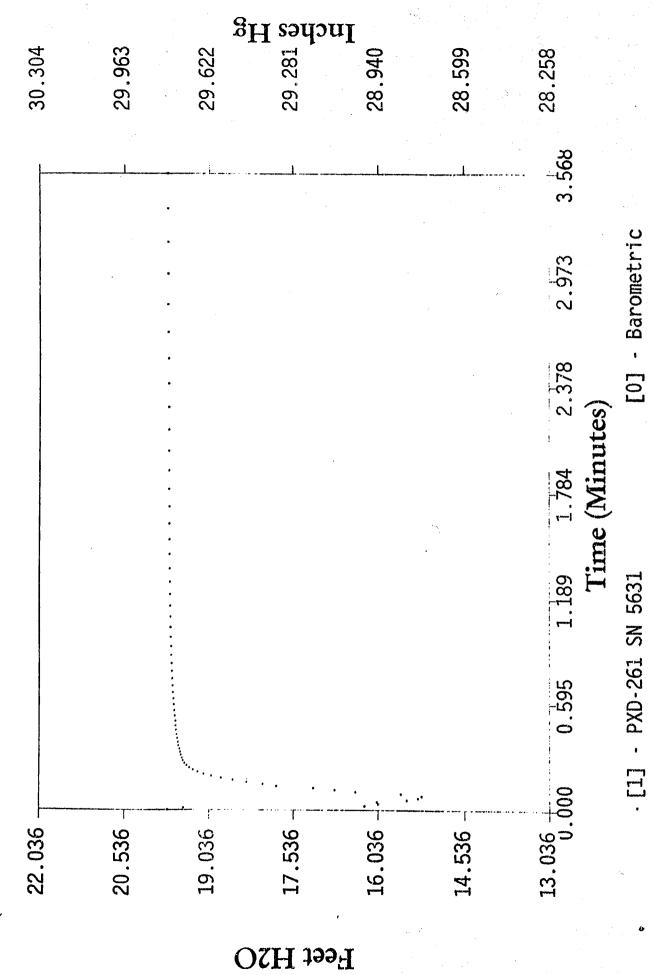
-1.55E+1 slope: intercept: 1.57E+0

Hydraulic Conductivity Bouwer-Rice: 7.35E-3 Hvorslev: 6.46E-3

### Recovery Data and Fit

time	У	weight	fitted
0.00	4.50	1.00	4.74
0.01	4.14	1.00	4.05
0.02	3.35	1.00	3.41
0.03	2.98	1.00	2.86
0.04	2.60	1.00	2.41
0.06	1.95	1.00	2.03
0.07	1.69	1.00	1.71
0.08	1.41	1.00	1.44

### 1S OUT



### Slug Test Results

Title: Slug out
Client: DMC
Job Number: 98-516
Well Number: MW-1 Deep

### Well Geometry

H: 300. Le: 12. Lw: 10.01 rc: .083 rw: 2.625

filter pack porosity: 0.0 effective radius: 8.33E-2

### Bouwer Rice Coefficients

Le/rw: 4.571 A: 1.721 B: .205 C: .738 ln(Re/rw): .71

### Hvorslev

F: 48.131

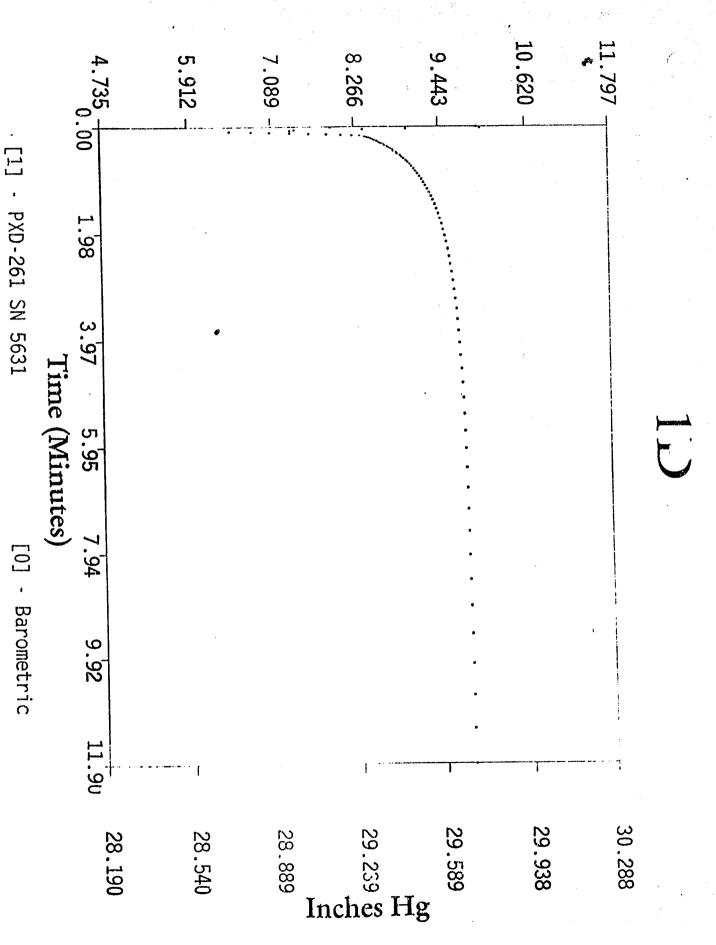
Least Squares Fit slope: -1.17E+0 intercept: 4.35E-1

Hydraulic Conductivity
Bouwer-Rice: 2.4E-4
Hvorslev: 5.3E-4

### Recovery Data and Fit

time	У	weight	fitted
0.00	1.56	1.00	1.54
0.01	1.53	1.00	1.53
0.02	1.51	1.00	1.51
0.03	1.48	1.00	1.49
0.04	1.47	1.00	1.47
0.06	1.44	1.00	1.45
0.07	1.43	1.00	1.43
0.08	1.39	1.00	1.40
0.10	1.38	1.00	1.38
0.11	1.36	1.00	1.36
0.13	1.33	1.00	1.33
0.14	1.30	1.00	1.31
0.16	1.28	1.00	1.28
0.18	1.25	1.00	1.25
0.20	1.23	1.00	1.22
0.22	1.20	1.00	1.19

Feet H2O



.

### Slug Test Results

Title: Slug in DMC Client: Job Number: 98-516

Well Number: MW-2 Shallow

Well Geometry

H: 38.17 12. Le: 38.17 Lw: .083 rc: 3. rw:

filter pack porosity: 0.0 effective radius: 8.33E-2

### Bouwer Rice Coefficients

Le/rw: 4. 1.705 A: B: .201 .71 C: 1.639

ln(Re/rw):

Hvorslev

52.228 F:

Least Squares Fit

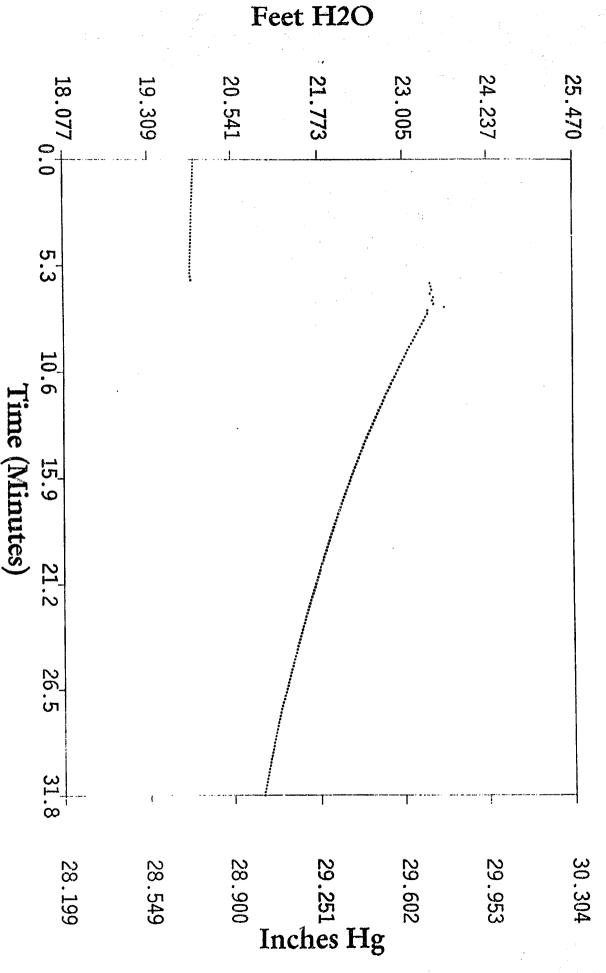
-4.8E-2 slope: intercept: 1.24E+0

Hydraulic Conductivity Bouwer-Rice: 2.28E-5 Hvorslev: 2.E-5

### Recovery Data and Fit

time	У	weight	fitted
0.00	3.45	1.00	3.45
0.17	3.42	1.00	3.42
0.33	3.40	1.00	3.39
0.5 <b>0</b>	3.37	1.00	3.37
0.67	3.34	1.00	3.34
0.83	3.31	1.00	3.31

# DMC MW-2 Shallow



- PXD-261 SN 5631

[0] - Barometric

### Slug Test Results

Title: Slug in
Client: DMC
Job Number: 98-516
Well Number: MW-2 Deep

### Well Geometry

H: 300. Le: 12. Lw: 31.94 rc: .083 rw: 2.625

filter pack porosity: 0.0 effective radius: 8.33E-2

### Bouwer Rice Coefficients

Le/rw: 4.571 A: 1.721 B: .205 C: .738 ln(Re/rw): .977

### Hvorslev

F: 48.191

Least Squares Fit

slope: -3.71E+1 intercept: 1.24E+0

Hydraulic Conductivity
Bouwer-Rice: 1.05E-2
Hyorslev: 1.68E-2

### Recovery Data and Fit

time	У	weight	fitted
0.00	3.43	1.00	3.31
0.01	2.19	1.00	2.28
0.02	1.48	1.00	1.50
0.03	1.02	1.00	0.99

Title: Slug out
Client: DMC
Job Number: 98-516
Well Number: MW-2 Deep

### Well Geometry

H: 300. Le: 12. Lw: 31.94 rc: .083 rw: 2.625

filter pack porosity: 0.0 effective radius: 8.33E-2

### Bouwer Rice Coefficients

Le/rw: 4.571 A: · 1.721 B: .205 C: .738 ln(Re/rw): · .977

### Hvorslev

F: 48.191

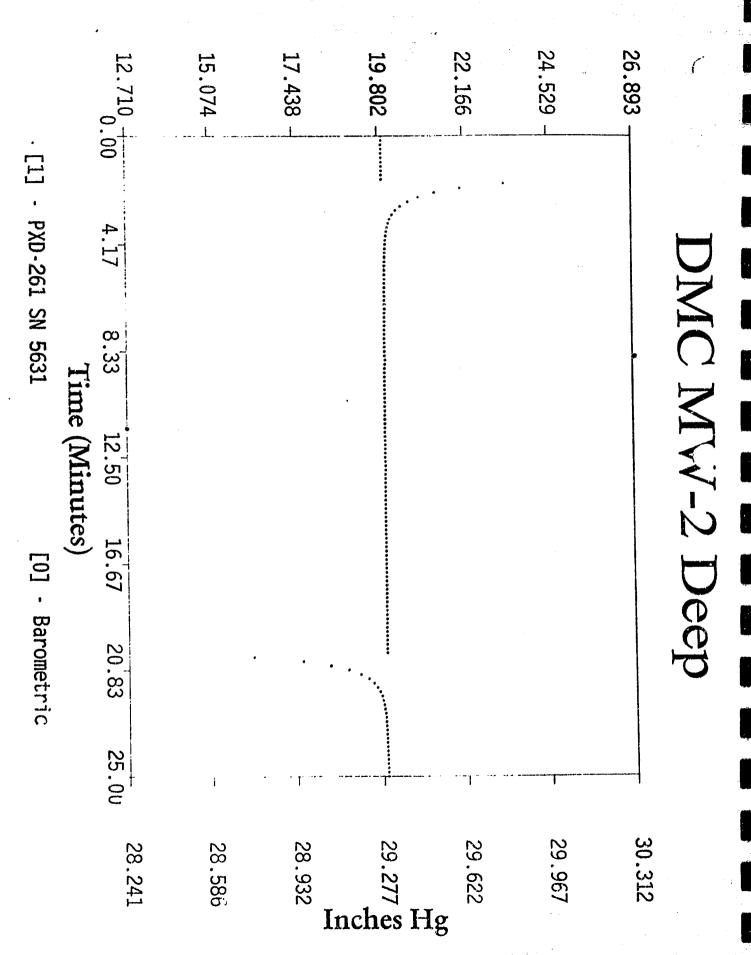
Least Squares Fit

**slope:** -3.84E+1 **intercept:** 1.31E+0

Hydraulic Conductivity
Bouwer-Rice: 1.08E-2
Hyorslev: 1.74E-2

У	weight	fitted
3.68	1.00	3.56
2.33	1.00	2.42
1.55	1.00	1.57
1.05	1.00	1.02
	2.33 1.55	3.68 1.00 2.33 1.00 1.55 1.00

Feet H2O



Title: Slug in Client: DMC Job Number: 98-516

Well Number: MW-3 Shallow

Well Geometry

H: 34.79 Le: 12. Lw: 34.79 rc: .083 rw: 3.

filter pack porosity: 0.0 effective radius: 8.33E-2

Bouwer Rice Coefficients

Le/rw: 4.
A: 1.705
B: .201
C: .71
ln(Re/rw): 1.597

III (Ne/ Lw) . . . .

Hvorslev

F: 52.228

Least Squares Fit

slope: -3.33E-1 intercept: 1.34E+0

Hydraulic Conductivity
Bouwer-Rice: 1.54E-4
Hvorslev: 1.39E-4

time	У	weight	fitted
0.00	3.90	1.00	3.83
0.17	3.61	1.00	3.62
0.33	3.39	1.00	3.43
0.50	3.20	1.00	3.24
0.67	3.04	1.00	3.07
0.83	2.90	1.00	2.90
1.00	2.79	1.00	2.74

Title: Slug out
Client: DMC
Job Number: 98-516

Well Number: MW-3 Shallow

### Well Geometry

H: 34.79
Le: 12.
Lw: 34.79
rc: .083
rw: 3.

filter pack porosity: 0.0 effective radius: 8.33E-2

### Bouwer Rice Coefficients

Le/rw: 4.
A: 1.705
B: .201
C: .71
ln(Re/rw): 1.597

Hvorslev

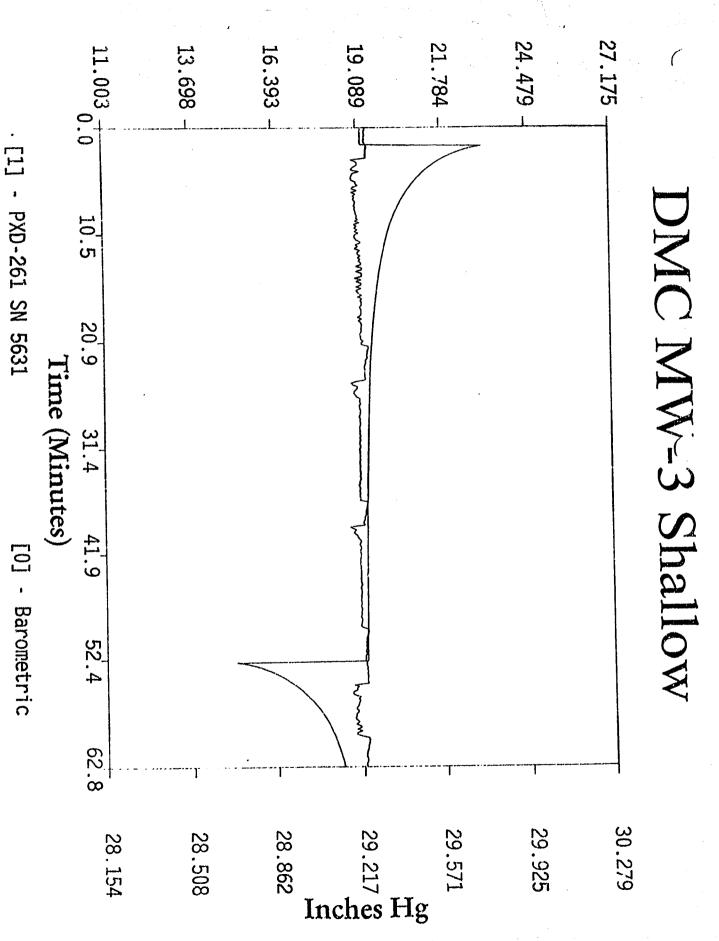
F: 52.228

Least Squares Fit slope: -3.35E-1 intercept: 1.41E+0

Hydraulic Conductivity
Bouwer-Rice: 1.55E-4
Hyorslev: 1.4E-4

time	V	weight	fitted
0.00	4.20	1.00	4.10
0.17	3,86	1.00	3.88
0.33	3.62	1.00	3.67
0.50	3,43	1.00	3.47
0.67	3.26	1.00	3.28
0.83	3.11	1.00	3.11
1.00	2.98	1.00	2.94

Feet H2O



Title: Slug in Client: DMC Job Number: 98-516 Well Number: MW-3 Deep

Well Geometry

H: 300. Le: 12. Lw: 35.9 rc: .083 rw: 2.625

filter pack porosity: 0.0 effective radius: 8.33E-2

Bouwer Rice Coefficients

Le/rw: 4.571 A: 1.721 B: .205 C: .738 ln(Re/rw): .997

Hvorslev

F: 48.191

Least Squares Fit slope: -2.12E+0 intercept: 9.47E-1

Hydraulic Conductivity
Bouwer-Rice: 6.1E-4
Hyorslev: 9.58E-4

time	v	weight	fitted
0.00	2.61	1.00	2.57
0.17	1.78	1.00	1.81
0.33	1.25	1.00	1.27
0.50	0.91	1.00	0.90

Title: Slug out
Client: DMC
Job Number: 98-516
Well Number: MW-3 Deep

Well Geometry

H: 300. Le: 12. Lw: 35.9 rc: .083 rw: 2.625

filter pack porosity: 0.0 effective radius: 8.33E-2

Bouwer Rice Coefficients

Le/rw: 4.571 A: 1.721 B: .205 C: .738 ln(Re/rw): .997

Hvorslev

F: 48.191

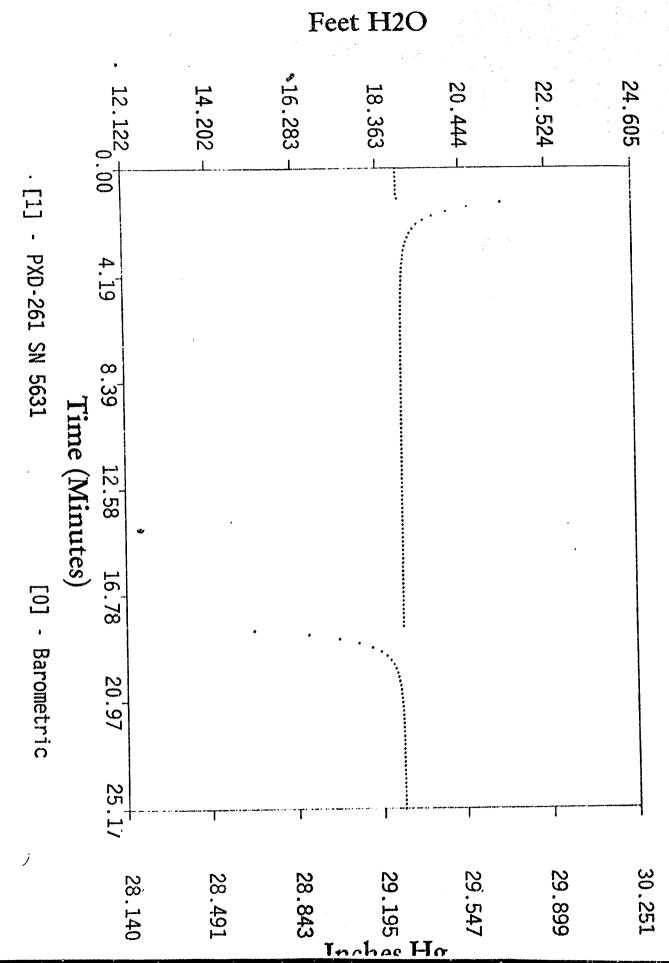
Least Squares Fit

slope: -2.42E+0 intercept: 1.28E+0

Hydraulic Conductivity
Bouwer-Rice: 6.98E-4
Hyorslev: 1.1E-3

time	У	weight	fitted
0.00	3.66	1.00	3.58
0.17	2.34	1.00	2.39
0.33	7ر'. 1	1.00	1.60
0.50	1.09	1.00	1.07

## DMC MW-3 Deep



3

Title: Slug in Client: DMC Job Number: 98-516

Well Number: MW-4 Shallow

### Well Geometry

H: 29.93 Le: 12. Lw: 29.93 rc: .083 rw: 2.125

filter pack porosity: 0.0 effective radius: 8.33E-2

### Bouwer Rice Coefficients

Le/rw: 5.647 A: 1.749 B: .211 C: .792 ln(Re/rw): 1.798

### Hvorslev

F: 42.813

Least Squares Fit

slope: -3.74E+0 intercept: 1.33E+0

Hydraulic Conductivity
Bouwer-Rice: 1.95E-3
Hyorslev: 1.91E-3

time	V	weight	fitted
0.01	3.57	1,00	3.64
0.17	2.06	1.00	2.03
0.33	1.12	1.00	1.09
0.50	0.57	1.00	0.58

Title: Slug out Client: DMC Job Number: 98-516

Well Number: MW-4 Shallow

### Well Geometry

H: 29.93 Le: 12. Lw: 29.93 rc: .083 rw: 2.125

filter pack porosity: 0.0 effective radius: 8.33E-2

### Bouwer Rice Coefficients

Le/rw: 5.647 A: 1.749 B: .211 C: .792 ln(Re/rw): 1.798

Hvorslev

F: 42.813

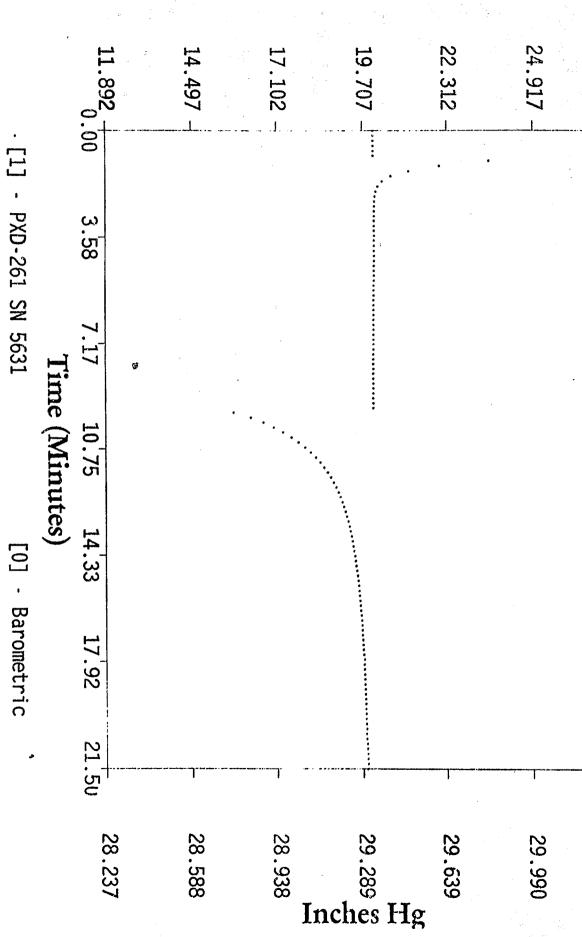
Least Squares Fit slope: -6.4E-1 intercept: 1.44E+0

Hydraulic Conductivity
Bouwer-Rice: 3.33E-4
Hvorslev: 3.26E-4

time	y	weight	fitted
0.01	4.26	1.00	4.18
0.17	3.75	1.00	3.79
0.33	3.36	1.00	3.40
0.50	3.04	1.00	3.06
0.67	2.75	1.00	2.75
0.83	2.50	1.00	2.47

Feet H2O 24.917 27.522 22.312 19.707 DMC MV-4 Shallow

30.340



Title: Slug in Client: DMC
Job Number: 98-516
Well Number: MW-4 Deep

### Well Geometry

H: 300. Le: 12. Lw: 26.03 rc: .083 rw: 2.625

filter pack porosity: 0.0 effective radius: 8.33E-2

### Bouwer Rice Coefficients

Le/rw: 4.571 A: 1.721 B: .205 C: .738 ln(Re/rw): .94

### Hvorslev

F: 48.191

Least Squar Fit slope: -3.E+0 intercept: 5.14E-1

Hydraulic Conductivity
Bouwer-Rice: 8.17E-4
Hyorslev: 1.36E-3

time	У	weight	fitted
0.01	2.52	1.00	2.44
0.17	1.48	1.00	1.53
0.33	0.90	1.00	0.93
0.50	0.58	1.00	0.56

Title: Slug out
Client: DMC
Job Number: 98-516
Well Number: MW-4 Deep

Well Geometry

H: 300. Le: 12. Lw: 26.03 rc: .083 rw: 2.625

filter pack porosity: 0.0 effective radius: 8.33E-2

Bouwer Rice Coefficients

Le/rw: 4.571 A: 1.721 B: .205 C: .738 ln(Re/rw): .94

Hvorslev

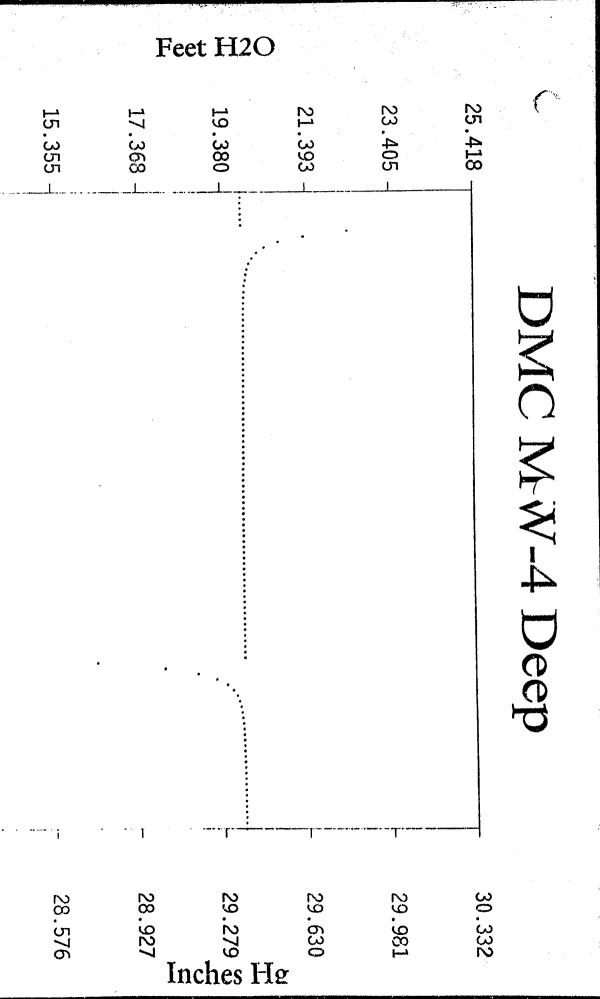
F: 48.191

Least Squares Fit

slope: -3.35E+0 intercept: 1.26E+0

Hydraulic Conductivity
Bouwer-Rice: 9.1E-4
Hvorslev: 1.52E-3

time	У	weight	fitted
0.01	3.54	1.00	3.40
0.17	1.93	1.00	2.01
0.33	1.12	1.00	1.15
0.50	0.68	1.00	0.66



13.343 + 0.00

3.19

6.39

9.58

12.78

15.97

19.17

28.225

Time (Minutes)

[0] - Barometric

[1] - PXD-261 SN 5631

Title:

Slug in

Client:

DMC

Job Number: 98-516

Well Number: MW-2 Mid

### Well Geometry

H:

39.03

Le:

3.5

Lw:

39.03

rc:

.083

rw:

2.125

filter pack porosity:

0.0

effective radius:

8.33E-2

### Bouwer Rice Coefficients

Le/rw:

1.647

A:

B: C:

ln(Re/rw):

2.646

### Hvorsley

F:

29.285

### Least Squares Fit

slope:

-3.17E-4

intercept:

1.86E+0

### Hydraulic Conductivity

Bouwer-Rice: 8.33E-7

Hvorslev:

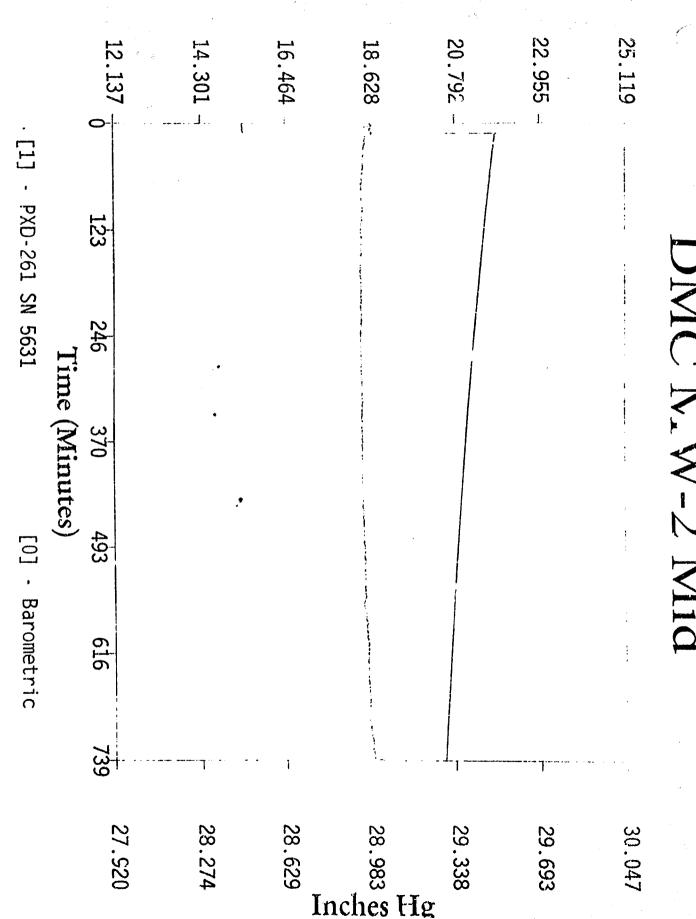
2.36E-7

time	у	weight	fitted
0.00	5.35	0.00	6.40
0.17	6.45	1.00	6.39
0.33	6.41	1.00	6.39
0.50	6.41	1.00	6.39
0.83	6.41	1.00	6.39
1.50	6.41	1.00	6.39

2.00	6.40	1.00	6.39
2.50	6.40	1.00	6.39
3.50	6.40	1.00	6.39
4.50	6.40	1.00	6.39
5.50	6.40	1.00	6.38
6.00	6.39	1.00	6.38
6.50	6.39	1.00	6.38
7.50	6.39	1.00	6.38
8.00	6.39	1.00	6.38
8.50	6.39	1.00	6.38
9.50	6.38	1.00	6.38
10.50	6.38	1.00	6.37
11.33	6.38	1.00	6.37
11.50	6.38	1.00	6.37
12.50	6.38	1.00	6.37
13.17	6.37	1.00	6.37
13.50	6.37	1.00	6.37
14.50	6.37	1.00	6.37
14.67	6.37	1.00	6.37
15.50	6.37	1.00	6.36
16.50	6.36	1.00	6.36
16.67	6.36	1.00	6.36
17.50	6.36	1.00	6.36
18.39	6.36	1.00	6.36
18.50	6.36	1.00	6.36
19.50	6.35	1.00	6.36
20.33	6.35	1.00	6.35
20.50	6.35	1.00	6.35
21.50	6.35	1.00	6.35
22.50	6.35	1.00	6.35
23.50	6.34	1.00	6.35
24.50	6.34	1.00	6.35
24.67	6.34	1.00	6.35
25.50	6.34	1.00	6.34
26.50	6.34	1.00	6.34
27.33	6.33	1.00	6.34
27.50	6.33	1.00	6.34
28.50	6.33	1.00	6.34
29.50	6.33	1.00	6.34
39.67	6.33	1.00	6.32
30.50	6.33	1.00	6.33
72.00	6.25	1.00	6.25
102.00	6.18	1.00	6.19
132.00	6.11	1.00	6.13

162.00	6.05	1.00	6.07
191.33	5.99	1.00	6.02
222.00	5.93	1.00	5.96
251.17	5.87	1.00	5.91
282.00	5.81	1.00	5.85
312.00	5.76	1.00	5.79
341.50	5.71	1.00	5.74
372.00	5.65	1.00	5.68
401.50	5.61	1.00	5.63
432.00	5.55	1.00	5.58
462.00	5.51	1.00	5.52
492.00	5.46	1.00	5.47
522.00	5.41	1.00	5.42
552.00	5.37	1.00	5.37
581.83	5.32	1.00	5.32
611.50	5.28	1.00	5.27
641.83	5.24	1.00	5.22
672.00	5.20	1.00	5.17
702.00	5.15	1.00	5.12
731.17	5.12	1.00	5.07

# DMC N.W-2 Mid



Feet H2O

Attachment 4

Summary Of Analytical Data

Compound	WH1-S	MW1-S	MW1-S	MW1-S	tEW1-S	MY1-S	MW1-S	EW1-S	MW1-S	MY1-S	MWI-S	MW2-S	1/2Y/2-\$	MW2-S	MW2-3
	5/7/98	?/2498	10/25/58	2/16/99	7/26/99	9/9/99	11/1/99	3/2/09	6/23/00	9/29/00	11/30/60	5/7/94	7/2496	10/26/98	7/25/99
Acetone .	<20	<20	<20		<50	≪0	<50					<20			<50
Acrolein	- 41	<10	<50		<50	<50	<50					<50 <5.0			<50 <50
Acrylontrile Benzera	<5.0 <1.0	<10 40	<10 40.5	40.5	<50 <5.0	<50	<50	≪5.0	<50	<5.0	<5.0	<1.0		40.5	90
Bromoterzene	<10	<1.0	<1.0	<1.0	40	<5.0	<5.0	4.0				<1.0			<5.0
Bromochloromethane	<1.0	<1.0	<1.0	<1.0	<5.0	<50	<50					<1.0			<5.0
Bromodichloromettære	<10	<1.0	<1.0	<1.0	<5.0	<5.0	<5.0					<1.0			<20
Bromoform	<13	920	₹20	<b>120</b>	<5.0	<5.0	<5.0					<1.0			<50
Bromomothene	<1.0	4.0	<4.0			<10.0	<10.0					<1.0 <1.0		ļ	<10.0
n-butyberzene sec-butyberzene	<1.0	<1.0	4.0 <1.0	<10		<10.0 <10.0	<10.0 <10.0					<1.0			<10.0
sert-buty/benzene	<1.0	<1.0	<1.0			<10.0	<10.0					<1.0			<10.0
Carbon disulfide	80	<1.0	<1.0		<10.0	<10.0	<10.0					<20			<10.0
Carbon tetrachionide	<1.0	<1.0	403	40.3	<5.0	<5.0	450					<10			45.0
Chloroconzene	<1.0	<1.0	<1.0			<5.0	<5.0					<10			<50
Chorodiconomethane	<1.0		<1.0			<5.0	<5.0					<1.0			<10.0
Chloroethane Chloroform	<1.0	40	4.0 41.0		<10 0 <5.0	<10.0 <5.0	<10.0 <5					<1.0			<50
Chloromethana	<1.0					<10.0	<10.0					<1.0			<10.0
6-chlorololuene	<10	<1.0	<1.0	<1.0	<5.0	<5.0	<5.0		ļ			<1.0	<del></del>		<5.0
p-chilorotoluene	<10				<5.0	<5.0	<50					<1.0			<50
1,2-dibramo-3-chiorapropene	<1.0											<1.0			
12-doromoethane	*1.0											<1.0	<del></del>	ļ	<b></b>
Disconomerane	<10				<0.0 <5.0	<5.0	<5.0 <5.0	<del></del>	<b> </b>		L	<1.0		<del> </del>	<u>- ব্</u>
1.2-dichlorubenzene 1.3-dichlorubenzene	<1.0	<10			G0 G0	<b>4.0</b>	<5.0		<del> </del>			<1.0	<del></del>	<b></b>	<50
n-dichlorobenzene	<10			<1.0	<5.0	450	<50		<del> </del>			<1,0			<5.0
Dichlorofluoromethane	<1.0												L		
Dichlorodifluoromet ana	<1.0				<10.0	<10.0	<10.0					<1.0			<10.0
12-dichloroethene	<10		<1.0		<5.0	<5.0	<5.0					<10			<0.0
1,1-dichloroethane	<1.0	<1.0	<1.0	<1.0	<50	<5.0	<50	<5.0 <5.0	+- <del></del>	<5.0 <5.0	<5.0 <5.0	<1.0 <1.0	<del></del>	<10	<5.0 <5.0
as-1,2-dationoffern	<1.0	11.0	<10			<5.0	<50		<del></del>			<1.0		<del> </del>	<50
rans-1,2-dichlorostrane	<1.0	<1.0	<1.0		<50	<b>-50</b>	<50					<1.0			<5.0
trans-1,4-dichloro-2-bulane	<1.0	<10	<10	<10	<10.0	<10.0	<10.0								<10.0
d-isopropylether	<10	<del></del>	<1.0		L										L
1,2-dichipropropane	<10		<1.0			<5.0	<50 <50		<b>_</b>			<1.0 <1.0			<50 <50
2.2-dichloropropene	<1.0	<10	410	<10	<5.0	450	<50		<del> </del>			<1.0	+	<del> </del>	30
1,1-dichioropropere	<1.0		<1.0			<50	<b>~50</b>		<b></b>			<1.0	<del></del>		<50
1,3-dichloropropene					<,0.6		<100								<10.0
cis-1,3-cichloropropene	40	<b></b>	<10	<del></del>	<5.0	<50	<50					<1.0	<del></del>		<50
trans 1,3 dichloropropene	- 40	<1.0	<1.0 <1.0		<5.0 <5.0	<50 <50	<50		<del> </del>			<1.0		<b></b>	<50 <50
Ethylcenzene Heptene		<1.0	·	<del></del>				·	<del> </del>			11.0	<del> </del>		130
2-hexanore	720		<10	<del>                                     </del>	<50	<5.0	<50		<del> </del>			70			<50
Heuarthorobutadiene	<10	<5.0	<5.0		<100	<10.0	<10.0					<1.0			<10.0
icdomethere	<10		<10	<del></del>			<10.0					<1.0	<del></del>		<10.0
Isopropylkenzene	-10		<1.0	ļ		<10.0 <10.0	<10.0 <10.0		<del> </del>			<1.0 <1.0		ļ	<10.0
p scpropytioluene Methylene chloride	<10	<1.0		<del></del>		30	<50		<del> </del>			<1.0		<b></b>	<10.0 <5.0
Hetryl etryl telone	4		<10		<100	<10.0	<100					<20	4	<b> </b>	<10.0
WIBE		<1.0	<10												
Metryl sictury kelone	ļ	<10	<10	1	ļ			L	ļ			ļ	ļ		
4-treiny-2-perlandra	- GO	L	- 40	<b> </b>	<50 <50	<5.0 <5.0	<50 <50	ļ	ļ			90			<50
Tecratry's charan Nacetturione	<1.0			ļ		130	150		<del> </del>			<1.0		ļ	<50
In sursylvanzere	<10	<10		4 <del></del>		<b>10.0</b>	<100				<del></del>	<1.0		<del> </del>	<10.0
Styrene	<10			<1.0		<50	<50					<1.0		I	<50
1,112-letrachtorcetrane	<10						<5.0					<1.0			<5.0
1.1.2.2 let/activ/cettiane	<10	<10	<10	<10		<50	<60	ļ	<u> </u>	<b></b> _	ļ	<1.0		<u> </u>	(50
Tetractéricellere	<10		<u> </u>			<50 <50	<5.0 <5.0	45 (	<50	<5.0	<50	<1.0	1	ļ <u></u> -	450
Toluera 12,3 that for chargera	<10	\$2 <b>40</b>				<5.0	र र र र	***	1			<1.0 <1.0		<1.0	<50 <50
1.2.4 trichorobe/usine	<10					40	<50	·	<del> </del> -				1	<del> </del>	<50
1,1 1-suctionoethere	<1 C		<10	<10	<50	<50	د30	·50	<50	<5.0	<50	<1.0	<1.0	<1.0	<del></del>
1,1,2-Victikroethare	<10	<10	<10			<50	<50								<50
Trichloroethene	<10	<10	<10		<50	<5U	<50		ļ			<10	<del></del>	<10	
Trichturofluoromethane	150	40	41.0			<10.7 <5.0	<10 0 <5 0		<del> </del>			<50	4	<b> </b>	<10.0
1.2.3 trickoruprovine 1.2.4 trinetnykarizene	<10 <10	<1.0 <1.0	<1.0 <1.0			*100	<10.0		<del> </del>			<1.0		<del> </del>	<5.0
1,3 5-transity/benzene	<10	رن در و	(10			<100	<10.0					<1.0		<del> </del>	<10.0
Veryl scelate	<50		<10	····	<100	<170	<100		<u> </u>			<50		<b> </b> -	₹10.0
Estyl exercia												<1.0	<del></del>		
Vinyl atkande	(10	<1.0	<1.0	<1.0	<100	<10.0	<10.0					<1.0			<10.0
2-ctskiroestryterryt ettier					ļ. <b></b>				<b> </b>			<10	4	ļ	ļ
Ethyl methacrylate o rylene	<10				<10.0	<100	<100					<b>110</b>	4	<del> </del>	<10.0
m and n sylene	410				450	<50	<50	<del>- ,</del> -		<del> </del>	<del></del>	41.0		<del> </del>	<50
				<30	<50	<5.0	150		·			<1.0		·	<5.0

Compound	NFWZ-8 9/9/99	HW2-8 11/1/99	MAZ-S 3/2/00	MY72-5 6/23/08	9/29/00	MW/2-S 11/30/60	924/2-\$ 11/30/06	MW3-S 5/7/98	2-EVAL 5/17/13	MW3-S 7/24/98	MW3-\$ 10/26/98	MW3-S 7/25/99	MW3-S 9/9/59	MW3-S 11/1/99	1/2/06
osone	<50	<50			1			<20				≪0	<50	<50	
crolein	<50	<b>&lt;</b> 50						<50	1			<50	<50	<\$0	
cryloniske	<50	<50						<5.0				<\$0	<50	<0	
Benzene	<5.0	<5.0	<5.0	<5.0	<20		<5.0	<1.0		40.5	<b>40.5</b>	<5.0	<5.0	<5.0	<
Bromobenzene	<5.0	<5.0	i					<1.0				⋖.0	<5.0	<b>~5.0</b>	
eromochloromethane	<50	<5.0						<1.0				<5.0	<5.0	60	
Bromodichioromethane	<5.0	<50			<del></del> 1			<1.0				≪0	<5.0	5.0	
							$\overline{\cdot}$	<1.0				<5.0	<50	40	
Bromoform	<50	<5.0										<10.0	<10.0	<10.0	
Bromomethane	<10.0	<10.0						<1.0							
n-busylbanzene	<100	<10.0						<1.0				<10.0	<10.0	<10.0	
sec bur/kenzene	<10.0	<10.0						<1.0				<10'0	<10.0	<10.0	
terl-buiy/benzene	<10.0	<10.0						<1.0				<10.0	<del></del>	<10.6	
Carbon disulfice	<10.0	<10.0						<b>7</b> 0				<10.0	¢10.0	<10.0	
Carbon tetrachionide	<b>450</b>	<5.0						<10				€0	<5.0	40	
Orloroberzene	<5.0	<5.0						<1.0				<5.0	<5.0	<50	
Chlorodibromomethene	<50							<1.0				<5.0	<5.0	40	
			<del> </del>					<1.0				<10.0	<10.0	<10.0	
Chicrosthane	<10.0	<10.0										40	€0	40	
Chlorolom	<50		ļ					<1.0							
Chioromethane	<10.0	<10.0						<1.0				<10.0	<100	<10.0	
o-chlorofoluone	<50	<5.0	Γ					<1.0				<5.0	≪.∂	€.0	
p-chlorotoluene	<5.0	<50										<5.0	<50	<50	
12-ditromo-3-chloropropene	<u>├</u>	<del></del>	<b></b>				·	<1.0							
	ļ	<del> </del>	<del> </del>	<b> </b>				<1.0		<del></del>					
1,2-dicromoeshane	<del></del>	-	<del> </del>		ļi					<del>  </del>		Æ A	<b></b>	<i>7</i> £ A	
Obromomethere	<50	<del></del>	<del></del>	ļ	<u></u> ,			<1.0		<del> </del>		<5.0	4.0	<50	
12-6:d/accompane	<5.0	<del></del>		ļ				<1.0	L	L		<5.0	<5.0	<b>45.0</b>	<u> </u>
1,3-dichlorobenzene	<5.0	<5.0	L	L				<1.0		<u></u>		<5.0	<5.0	<5.0	L
o dichinobergane	<5.0	<5.0						<10				<5.0	<50	<50	L
Dichlorofluoromethane		1	<del> </del>	l		·········							_		
Dichlorud@usrome@iene	<100	<10.0	<del> </del>	<b></b>			<b></b>	<1.0				<10.0	<10.0	<10.0	· · · · · · · · · · · · · · · · · · ·
				<del></del>	<del></del>	<del></del>		<1.0				<5.0	<5.0	<5.0	
1,2-dichloroethane	<5.0				<u> </u>					<del>  </del>	4.0	<b>40</b>		45.0	$\vdash$
1,1-dichlorpethane	<5.0	<5.0			⋖.0		<50	2.9	2.0	ļ	4.0				
1,1-dishlorosthene	<b>45</b> .,	<5.0	<5.0	<50	<50		<50	<1.0		<2.0		<5.0	<5.0	<b>450</b>	
as-1,2-diaticroethene	<50	<b>450</b>						<1.0	1	]		<5.0	<.0	⋖5.0	l
vars-1,2-dichlorosthers	<5.0	<5.0						<10				<5.0	<50	<5.0	
	<10.0											<10.0	<10.0	<100	
rans-1,4-dichloro-2-bulana	10.0	10.0	<del> </del>		ļ		<del> </del>			/ <del></del>		—— <del>"</del>			
d isopropylether	<u> </u>		<b>├</b> ──		ļ		ļ		<u> </u>	<del> </del> -		<del></del>			
1,2-dichloropropene	<50		+			ļ	<b> </b>	<10	<u> </u>			<50	<50	<5.0	ļ
1,3 dichioropropana	<50	<50	1				<u> </u>	<1.0				<50	<50	<5.0	
2.2-dichioropropane	<50	<50	·		l	·		<10				<50	<50	<50	<u> </u>
1,1-dichloropropens	<5.0	<50		T				[				<5.0	<b>G</b> 0	≪50	
1,3 dichloropropene	<10.0			<u> </u>				i				<10.0	<10.0	<10.0	
<del></del>	<5.0	+		<del></del>				<1.0	7.			<50	<b>45</b> 0	450	
cus-1,3-dichloropropene	<del></del>	<del></del>	<del></del>	<del> </del>			<del> </del>					40	·	<5.0	
vens-1,3-dictioropropene	<5.0	<del></del>		<b></b>			ļ	<1.0					<del></del> -		
Ethykenzene	<50	<50	1	1	<u> </u>			<1.0			·	<5.0	<b>450</b>	<50	<u> </u>
Hectane	i	l	l	L	l		<u> </u>								<u> </u>
2-hezarone	<b>150</b>	<50			]			<20		[		<50	450	< 0	
Heuschkrobutediene	<10.0	<10.0					i	<10				<10.0	<100	<10.0	1
kdometrare	<10.0	<100		<del> </del>		·		<1.0				<10.0	<10.0	<10.0	
	<17,0	<del></del>	+					<1.0	<del></del>	<del> </del>		<10.0		<100	<del> </del>
lvaropyturare		<del></del>	<del></del>	<del> </del>		<del> </del>			<del></del>			<10.0		<10.0	<del> </del>
p-reopropy/ficilitiene	4100	<10.0	<del></del>	<del> </del>	<del> </del>	<b> </b>	<del></del>	<10			<del></del>	<del></del>			ļ
Metrylana chkrida	(50	·	·	ļ	ļ	<b></b>	<b></b>	<1.0	<del></del>	ļ		<50		<50	
Mothyl ethyl ketone	<10.0	<100				<u> </u>		20	L	<u> </u>		<10.0	<100	<10.0	
WTEE	T	I	1		l	L			L			L	L	L	L
V etnyl rackutyl ketone	1	T	Γ		<u> </u>	Γ	T	1	T	[			1	I	Ι
4 matryl 2 pantamone	<5.0	<50	1	T			T	<20				<5.0	<50	<0	1
				<del> </del>	<del> </del>	<del> </del>	<del> </del>	<50	<del></del>	t		<50	<del></del>		
Tetrahydroluran	<del> `2°</del>	1	1	<del> </del>	<del> </del>	<del> </del>	<del> </del>	<10		<del> </del>		<del> </del>	<del> </del>	<del> </del>	!
liscetta ene	<b> </b>	ļ	<del> </del> -	<del> </del>	ļ		<del> </del>			<b> </b>			<del> </del>	<del></del>	<del> </del>
n propytuenzone	<10.0			<del> </del>	ļ		ļ	<1.0	+	<b> </b>		<10.0			
Streen	<30						L	<1.0		<b></b>		40	<del></del>	<50	
1 1 1,2 tetrarticroetrane	<b>~50</b>	30					L	<10				40	<b>40</b>	<50	
1,1.2.2-letrachiorcethare	30	\$	<del></del>	Ι	1		1	<1.0				ব			+
	₹ 0			<del> </del>	<del> </del>	<del></del>		×1.0		T		40			
Tevachkircethene				<del> </del>		<del> </del>	<50	<1.0	<del></del>	<1.0	<10	<del></del>		+	
Total e	<50	<del></del>		<5.0	<50	ļ		<del></del>	<del></del>	****		<del>4</del>			
12.3 trichiorobenzene	<5.0	<del></del>	<del></del>	<u> </u>	ļ		<b></b>	<10		<b> </b>		্ ব	+		
1,2 4 trichkeubenzene	<50			1	!	<u></u>	<u> </u>	<1.0				<50	4		
1 1 1 Inchlorce/hane	<50	<50	(5)	<50	<50	L	<50	15	12.0	12.7	11.4	<5(	6.80	6.4	
1.1.2 trictionoethene	<50	·	1	!	I		I	<10		1		30	<50	450	
Trichlorcethere	<50	<u> </u>	<del></del>	F=		·		<1.0		<10	<1.0		+		
	1100	<del></del>	<del></del>	<del> </del>	<del> </del>	<del> </del>	<del> </del>	75.0	<del></del>	<u>`</u>	<del>-</del>	<100	+		<del></del>
Inchlorofucromethane	<del></del>	<del></del>		·	<del> </del>	<del> </del>	<del></del>		<del></del>	<del> </del> -			<del></del>	<del></del>	
1,2,3 trichloropropane	<50	<del></del>		<b> </b>	ļ	ļ	<u> </u>	<1.0		<b></b>		<50			-
1,2,4-tnmstrytenzene	<100	<10.0	<u></u>	L		ļ. <u></u>		<1.0	<del></del>			<100	<100	<10.0	1
1.3.5 unietrykerzere	<10.0	<100	T					<10				<100	<10.0	<100	
	<b>4100</b>			t	<del> </del>	1	1	<50	·	T	·	<10 (	<del></del>	4	+
Virigi acetate	*100	*100		<del> </del> -		<del> </del>	<del> </del>		<del></del>	<del> </del>	<del> </del>	<del> </del>	1	1	<del> </del>
try scalale	<u> </u>		<b></b>	ļ		ļ	ļ	<1.0		<b> </b>	<b> </b>	ļ	ļ	<b> </b>	<del> </del>
/inyl chiande	(100	<10.0	<b></b>	<b></b>	ļ	ļ	<b></b>	<1.0		ļ	.,	<10.0	<10.0	<10.0	4
-crici cethylwnyl ett er	1	I	L	L	L	L	L	<1.0	L	L	L		L	L	
	I	l	l	l				<5	1	1				1	
thyl methachisia		L	<del></del>			†- <del></del>	<del> </del>		<del></del>	<del></del>	<del></del>	·		1	·
thyl methecrylale	£10.0	<10.0	i	1	i i	1		<b>1 (10</b>	1			<100	)] <10.0	(10.0	
thyl methacrylala xylaria and p xylana	<10.0 <5.0	<10.0 <5.0				<b> </b>	ļ	<1.0 <1.0				<100			

Compound	MW3-S 11/1/59	MW3-S 3/2/00	MW3-S 6/23/00	MW3-S 8/29/00	MW3-S 11/30/00	MW4-S 5/7/98	MW4-S 7/24/98	MW4-S 10/26/98	MW4-S 7/26/99	MW4-S 9/9/99	MW4-S 11/1/99	MW4-S 3/2/00	MW4-S 6/23/00	MW4-8 9/29/00
Acetone	117/99 <50	3/2/00	WZ3/JV	DIEZKAN	11130100	<20	114-170	10120130	<50	<50	<50			
Acrolein	450					<50			50	<50	<30			
Acrylonitile	<50					<5.0			<50	<50	<50			
Benzere	<5.0	<5.0	<5.0	<5.0	<5.0	<1.0	<0.5	<0.5	<5.0	<5.0	<5.0	<5.0	<5.0	₹5.0
Bromobenzane	<5.0					<1.0			<5.0	<5.0	<b>45.0</b>			
Bromochloromethane	<5.0					<1.0			<5.0	<5.0	<5.0			
Bromodichloromethane	<5.0					<1.0			<5.0	<5.0	<5.0			
Bromoform	<5.0					<1.0			<5.0	<5.0	<5.0			
Promometrane	<10.0					<1.0			<10.0	<10.0	<100			
n-butybenzene	<10.0					<1.0			<10.0	<10.0	<10.0			
sec-buty/benzene	<10.0					<1.0			<10.0	<10.0	<10.0			
tert-butyfbenzene	<10.0					<1.0			<10.0	<10.0	<10.0			
Carbon disutfide	<10.0					<20			<10.0	<10.0	<10.0			
Carbon tetrachloride	<5.0					<1.0			<5.0	<5.0	<5.0			
Chiorobenzene	<5.0					<1.0			<5.0	<5.0	<5.0			
Chlorodibromomethane	<5.0					<1.0			<5.0	<5.0	<5.0			
Chloroethane	<10.0					<1.0			<10.0	<10.0	<10.0			
Chloroform	<b>45.0</b>					<1.0			<5.0	<5.0	<5.0			
Chloromethane	<10.0		L			<1.0			<19.0	<10.0	<10.0			
o-chiorotoiuene	<5.0					<10			<5.0	<5.0	<5.0			
p-chicrotoluene	<5,0	L				<1.0			<5.0	<5.0	<5.0			
1,2-dibromo-3-chloropropane						<1.0							L	
1,2-dibromoethane						<b>&lt;1.0</b>			<b> </b>					<u> </u>
Dibromomethane	<5.0		ļ	ļ		<1.0			<5.0	<5.0	<5.0		<u> </u>	
1,2-dichlorobenzene	<5.0				ļ.,	<1.0	<b> </b>		<5.0	<5.0	c5.0		ļ	
1,3-dichlorobenzene	₹5.0					<1.0	<u> </u>		<5.0	<5.0	<5.0		ļ	
p-dichlorobenzene	<5.0	L		L	L	<1.0	ļ		<5.0	<5.0	<5.0		<u> </u>	
Dichlorofluoromethane			<u> </u>				<u></u>							
Dichlorodifluoromethane	<10.0		<u></u>			<1.0	<b></b>		<10.0	<10.0	<10.0		ļ	
1,2-dichloroethane	<5.0			<u> </u>		<1.0			<5.0	<5.0	<5.0			
1,1-dichkroethane	<5.0				<5.0	<1.0		<1.0		<5.0	₹5.0	<5.0	<5.0	<5.0
1,1 schloroethene	<5.0		<5.0	<5.0	<5.0		<2.0		<5.0	<5.0	<5.0	<5.0	<5.0	<5.0
r.s-1,2-dichloroethene	<5.0				L	<1.0			15.0	<5.0	<5.0			
trans-1,2-sichloroethene	<5.0		ļ	ļ		<1.0	Ļ		<5.0	<5.0	<5.0		ļ	ļ
rans-1,4-dichloro-2-butane	<10.0	<u> </u>	<u> </u>	<b></b>	ļ	<u> </u>	ļ		<10.0	<10.0	<10.0	<b> </b>	<del> </del>	
di-Isopropylether		ļ		<u> </u>	ļ	ļ								ļ
1,2-dichloropropane	<5.0		<del> </del>	ļ		<1.0	ļ		<5.0	<5.0	<5.0		<b> </b>	
1,3-dichloropropane	<5.0		ļ	ļ		<1.0			<5.0	<5.0	<b>45.0</b>			
2,2-dichloropropane	<5.0		ļ	<b></b>		<1.0			<5.0	<5.0	<5.0		ļ	<b> </b>
1,1-dichloropropene	<5.0		ļ	<del> </del>		<1.0	<b> </b>		<5.0	<5.0	<5.0		<u> </u>	ļ
1,3-dichloropropene	<10.0			<b> </b>		<del> </del>	<del> </del>		<10.0	<10.0	<10.0		<b>}</b>	
cis-1,3-dichloropropene	<5.0		<b> </b>	<b></b>		<1.0		<u> </u>	<5.0	<5.0	<b>&lt;</b> 5.0		<del> </del>	
trans-1,3-richkropropene	<5.0		<del> </del>	<del> </del>	<del> </del> -	<1.0	1	<del></del>	<5.0	<5.0 <5.0	<5.0 <5.0		<b> </b>	
Ethylbenzane	<5.0	<del>' </del>	<b></b>	<del> </del>		<1.0	<del> </del>	<del> </del> -	<5.0	45.0	45.0	<del> </del>	<del> </del>	ļ
Heptane	<5.0	<del> </del>	<b>}</b>	<del> </del>	<del> </del>	<20	<del> </del>	<del> </del>	<5.0	<5.0	<5.0		<del> </del>	<del> </del>
2-hexanone	<10.0		<del> </del>	ļ		<1.0		ļ	<b>₹10.0</b>		<10.0		<del> </del>	<del> </del>
Hexachlorobutadiene lodomethane	<10.0		<del> </del> -	<del> </del>	<del> </del>	<1.0			<10.0		<10.0		<del> </del>	
Isopropylbenzene	<10.0		<del> </del>	<del> </del>	<del> </del>	<1.0		<del> </del>	<10.0				<del> </del>	<del> </del>
p-isop-opytokiene	<10.0		L	<del> </del>	<del>├─</del> ─	<1.0		<del></del>	<10.0		<10.0		<del> </del>	<del> </del>
Methylene chloride	<5.0		<del></del>	<del> </del>	<del> </del>	<10			<5.0				<del> </del>	
Metryl ethyl kalone	<10.0		<del> </del>	<del> </del>	<del> </del>	<20		<del> </del>	<10.0				<del>                                     </del>	<del>                                     </del>
MIBE	<del></del>	1	<del> </del>	<del>                                     </del>	<del> </del>	1	<del>                                     </del>	<del>                                     </del>	1	1	1	<del> </del>	<del>                                     </del>	<del>                                     </del>
Methyl isobutyl ketone	1	1	<del>                                     </del>	<del> </del>	<del> </del>	<20	1	<del>                                     </del>	<del> </del>	<del>                                     </del>	<del> </del>	<del>                                     </del>	<del>                                     </del>	<del>                                     </del>
4-metryl-2-pentanone	<5.0	<u> </u>	T	<del>                                     </del>	<del> </del>	<20		<u> </u>	<5.0	<5.0	<5.0	1	t	
Tetrahydrofuran	<5.0		Ī		1	<50	1	<u> </u>	<5.0	<5.0	<5.0		T	7
Naphthalene		L	I''			<1.0		L				<u></u>	Ι	I
n-propy(bertzene	<10.0		1	]	I	<1.0			<10.0	<10.0	<10.0		T	<u> </u>
Styrene	<5.0			1	<u> </u>	<1.0			<5.0			1	<u> </u>	
1,1,1,2-letrachloroethane	<5.0				I	<1.0			<5.0		<5.0		<u> </u>	
1,1.2,2-tetrachioroethane	<5.0					<1.0			<5.0				I	I
Tetrachloroethene	<5.0			1		<1.0			<5.0				T	
Toluera	<5.0		<5.	0 <5.0	<5.0	<1.0	<1.0	<1.0	<5.0	<5.0	<5.0		0 <5.	<5.0
1,2,3-trichlorobenzene	<5.0		L			<1.0			<5.0					
1,2,4-trichlorobenzene	<5.0					<1.0			<5.0					
1,1,1-trichloroethane	6.4	<5.0	<5	0 <5.0	<5.	<1.0	<1.0	<1.0	0 <5.0	<5.0	<5.6	<5.	u <5	0 ∢5.0
1,1,2-trichloroothane	<5.0					<1.0			<5.0				$\Box$	
Trichloroethene	<5.0					<1.0		<1.0						
1'nchlorofluoromethane	<10.0			1		<5.0			<10.0					
1,2,3-trichioropropane	<5.0					<1.0			<5,0			٥	<u> </u>	I
1,2,4-trimothythenzene	<10.0		L			<1.0			<10.0			)		J
1,3,5-trimethythenzene	<10.0		I	L		<1.0		L	<10.0	<10.0	<10.	0	I	T
Vinyi acetate	<10.0		I	I		<5.0			<10.0	<10.0	<10.	0	1	
Ethyl acotate	T	1		T	L	<1.0		I	<u> </u>		<u> </u>		1	T
Virryl chloride	<10.0		1	1	T	<1.0		1	<10.0	<10.0	<10.	0	1	1
2-chloroethyMnyl ether	<b>1</b>	1	1	1	T	41.0		1	1	1	1	1	1	1
Ethyl methacrylala	1	1	<del> </del>	1	1	4		1	1	1	1	1	1	1
)	<10.0	1	1	1	T	<1.0		1	<10.0	<10.0	<10.	0	1	1
In-XYIGNO								<del></del>					-	+
n-xylene m and p xyleng	<5.0			i i		<1.0	)	Į.	<5.0	) <5.0	) <6.	이	l .	

	MW4-S 11/30/00	MW5-S 7/26/99	MW5-8 9/9/99	MW5-S 11/1/99	MW5-S 3/2/00	MW5-S 6/23/00	MW5-S 9/29/00	MW5-S 11/30/00	MW6-S 7/26/99	MW6-S 9/9/99	MW6-8 11/1/99	1/2/00	14 <b>4</b> 6-5 5/23/00	MW6-S 9/29/00
Acatone	-	<50	₹50	<50					<50	ر50	<50			
Acrolein	<del>  </del>	<50	<50	<50					<50	<50	<50			
Acrylonitrile		<50	<50	<50					<50	<50	<50			
Benzene	<5.0	<5.0	<5.0	<5.0	<5.0	<5.0	<5.0	<5.0	<5.0	<5.0	<5.0	<5.0	<5.0	<5.0
Bromobenzene		<5.0	<5.0	<b>45.0</b>					<5.0	<5.0	<5.0			
Bromochloromethane	1	<5.0	<5.0	<5.0					<5.0	<5.0	<5.0			
Bromodichioromethane	1	<5.0	<5.0	15.0					<5.0	<5.0	<5.0			
Bromoform	<del>  </del>	<5.0	<5.0	<5.0					<5.0	<5.0	<5.0			
Bromomethane	1	<10.0	<10.0	<10.0					<10.0	<10.0	<10.0			
n-butytbenzene	1	<10.0	<10.C	<10.0					<10.0	<10.0	<10.0			
sec-butythenzene	<del>  </del>	<10.0	<10.0	<10.0					<10.0	<10.0	<10.0			
	<del> </del>	<10.0	<10.0	<10.0		<del>                                     </del>			<10.0	<10.0	<10.0			
lext-buly/benzone	<del> </del>		<10.0	<10.0		<del> </del>			<10.0	<10.0	<10.0			
Carbun disulfide	<b>↓</b>	<10.0								<5.0	<5.0			
Carbon tetrachloride		<5.0	<5.0	<5.0					<5.0		<5.0 <5.0	<del></del>		
Chlorobenzene	.	<5.0	<5.0	<5.0					<5.0	<5.0		<del></del>		
Chlorodibrornomethane		<5.0	<5.0	<5.0					<5.0	<5.0	<5.0			
Chloroethanu		<10.0	<10.0	<10.0					<10.0	<10.0	<10.0			
Chloroform		<5.0	<5.0	<5.0					<5.0	<50	<5.0			
Chloromethane	1	<10.0	<10.0	<10.0					<10.0	<10.0	<10.0			
o-chlorotokiene	1	<5.0	<5.0	<5.0					<5.0	<5.0	<\$.0			
p-chlorotolusne	1	<b>&lt;5.0</b>	<del>3.0</del>	<5.0	<b> </b>				<5.0	<5.0	<5.0			
<del></del>	+	****	- ~··		<b> </b>	<del> </del>			i		<del>-</del>			***************************************
1,2-dibromo-3-chioropropane	<del> </del>		<del> </del>		<del> </del>	<del> </del>				<b></b>	<del></del>			<del></del>
1,2-dibrornosthane	<del> </del>		<u> </u>		<del> </del>	<del> </del>			-50	<5.0	<5.0	<del> </del>		
Dibromomethane		<5.0	<5.0	<5.0	<b> </b>	<b> </b>			<5.0		<5.0	<del> </del>		
1,2-dichlorobenzene	<b></b>	<5.0	<5.0	<5.0		<b> </b>		. <u></u>	<5.0	<5.0		<del> </del>		<b></b>
1,3-dichlorobenzene	1	<5.0	<5.0	<5.0	<b></b>	<b> </b>			<5.0	<5.0	<5.0	<del> </del> -		
p-dichlorobenzene		<5.0	<5.0	<5.0					<5.0	<5.0	<5.0	<b> </b>		<u> </u>
Dichlorefluoromethane				L						L	<u> </u>	<b></b>		ļ
Dich(cro@fluoromethane		<10.0	<10.0	<10.0					<10.0	<10.9	<10.0			ļ
1,2-dichloroethane		<5.0	<5.0	<5.0					<5.0	<5.0	<5.0			
1,1-dichioroethane	<5.0	<5.0			<5.0	<5.0	<5.0	<5.0		<5.0	<5.0	<5.0	<5.0	<5.0
1,1-dichloroethene	<5.0	<5.0			<5.0		<5.0	<5.0	≺5.0	<5.0	<5.0	<5.0	<5.0	<5.0
cis-1,2-dichioroethene		<5.0				1			<5.0	<5.0	<5.0			
	<del></del>	45.0			<del> </del>	<del> </del>			<5.0		<5.0	<del></del>		
trans-1,2-dichloroethene						<del> </del>	ļ		<10.0					
trans-1,4-dichloro-2-butene	<b>_</b>	<10.0	<10.0	<10.0	<b> </b>	<del> </del>	ļ		\$10.0	10.0	10.0	<del> </del>		<del> </del> -
di-isopropylether			<u></u>						L	<u> </u>	<del> </del> _			<b></b>
1,2-dichloropropane		<5.0				1	<u> </u>		<5.0					<b></b> .
1,3-dichloropropane	1	<5.0						ļ	<5.0		<5.0			
2,2-dichloropropane		<5.0	<5.0	<5.0				<u> </u>	4.0	<5.0				
1,1-dict/oropropene		- 5.0	<5.0	<5.0	1			·	<5.0	<5.0	<5.0	L		
1,3-dichlorapropene	1	<10.0	<10.0	<10.0					<10.0	<10.0	<10.0			
cls-1,3-dichlatopropene	<del>-</del>	<5.0	<5.0	<5.0		1	1		<5.0	<5.0	<5.0			
trans-1,3-dichi vopropone	<del></del>	<5.0	<5.0	<5.0	1	<del></del>			<5.0	<5.0	<5.0			
Ethythenzene	<del></del>	<5.0				<del></del>		<u> </u>	<5.0	<5.0	<5.0			
Heptane	<del>- </del>	<del> </del>	<del> </del>		<del> </del>	·		<del></del>	† <u>-</u> -	<del></del>	<del>                                     </del>			
		<5.0	₹.,	<5.0	<del> </del>	<del> </del> -	<del> </del>		<5.0	<5.0	<5.0	<del> </del>		<del> </del>
2-hexanone		<10.0	<del></del>			<del> </del>	<del>}</del>	<del></del>	<10.0				<del> </del>	<del> </del>
Hexachlorobutariene			4			<del> </del>							<del></del>	
lodomethane		<10.0			<del></del>	<del> </del> -	<del> </del>	ļ	<10.0				ļ	<del> </del>
Isopropyfbenzene		<10.0	<del></del>				<del> </del>	ļ	<10.0			4	Ļ	<del></del>
p-isopropy totuene		<10.0	<del></del>			<u> </u>	L		<10.0	<del></del>				ļ
Methylene chloride		<5.0			4	<u> </u>	L		<5.0	·			<u> </u>	<b> </b>
Methyl ethyl kelone		<10.0	<10.0	<10.0	1		ļ		<10.0	<10.0	<10.0	1	ļ	<u></u>
MIBE			L				L						L	
Methyl isobutyl ketone		<u> </u>	1		I					L			L	
4-methyl-2-pentanone	T	<5.0	<5.0	<5.0	)]	T	I		<5.0	<5.0	<5.0	0	I	I
Totrahydrofuran	+	<5.0				1	1	I	<5.0				[	T
Naphthalene	<del></del>	<del>                                     </del>	† <del></del>	<del>                                     </del>	1	<del> </del>	<del> </del>	<del></del>	1	1	1	1	<del> </del>	1
	<del></del>	<10.0	<10.0	<10.1	<del>\ </del>	<del> </del>	<del> </del>	1	<10.0	<10.0	<10.0	1	<del>                                     </del>	<del> </del>
n-propylbenzene	<del></del> ·					+	<del> </del>	<del> </del>	<5.0				<del> </del>	<del> </del>
Styrese		45.0				+	<del> </del>	<del> </del>						1
1,1,1,2-letrachloroethane	4	<5.0				+	<del> </del>	<b>}</b>	<5.0				<del> </del>	<del> </del>
1,1,2,2-letrachloroethane	<b></b>	<5.0				<del> </del>	<b> </b>	<del> </del>	<5.0				ļ	<del> </del>
Tetrachiorcethene		<5.0				<del></del>	<b></b>		<5.0				<u> </u>	<b></b>
Toluene	<5.0					.0 <5.0	<5.0	<5.0					<5.0	0 <5
1,2,3-bichlorobenzene		<5.0							<5.0				<u></u>	
1,2,4-trichiorobenzene	T	<5.0		45.0					<\$.0	<5.0	< 5.	0		
1,1,1-trichloroethane	<5.0					.0 <5.0	<5.0	<5.0					<5.	0 <5
1,1,2-frichloroethane	- <del></del>	<5.0				1	1	1	<5.0				<del> </del>	1
	<del> </del> -	<5.0				1	1	1	48.0				<del> </del>	<del> </del>
Trichloroetherie	<del></del>	<10.0	·			+	<del> </del>	-	<10.0				<del> </del>	+
Trichlorofluoromelhane			4			<del> </del>	<del> </del>	<del> </del>					<del> </del>	<del> </del>
1,2,3-trichloropropane		<5.0				<del> </del>	<del> </del> -	<del> </del>	<5.0				<del> </del>	<del> </del>
1,2,4-timethylbenzene		<10.0				<del> </del>	<b>↓</b>	ļ	<10.0				<del> </del>	Į
1,3,5-trimethy/benzene		<10.0					<b></b>	<u> </u>	<10.0				<b> </b>	1
Vinyl acetate		<10.0	<10.0	<10.0	<u></u>		<u></u>	L	<10.1	<10.	0 <10.	0	J	
Ethyl acetate		1	1	1				1		1	1	.1	1	
	-	<10.0	<10.0	<10.0	1	1	1	T	£10.0	<10.	o <10.	0		<del> </del>
		1	4	1	<del> </del>	<del> </del>	<del></del>	<del> </del>	1	1	1		+	<del> </del>
Vinyt chloride	<del></del>	1	1	1										
Vinyt chloride 2-chloroethylvinyt ether		<del> </del>	<u> </u>	<del> </del>	<del> </del>	<del> </del>	<del> </del>	<del> </del>	<del> </del>	<del> </del>	<del> </del>	<del> </del>	<del> </del>	+
Vinyt chloride 2-chloroethylvinyt ether "ryt methacrytete			<u></u>										ļ	<del> </del>
Vinyt chloride 2-chloroethylvinyt ether		<10.0							<10.0 <5.0					

Compound	MW6-S	UW7-8	MW7-S	MW7-S	kW7-8	MW7-S	MW7-S	MW7-S	MW8-8	MW8-S	MW8-S	MW8-8	MW8-9	MW8-S
Quiposiio	11/30/90	7/26/99	9/9/99	11/1/99	1/2/00	6/23/00	9/29/00	11/30/00	7/20/99	9/9/99	11/1/99	3/2/00	8/23/00	9/29/00
Acelone		<50	<50	<50					<50	<50	<50			
Acrolein		<50	<50	<50					<50	<50	<50			
Acrylonitrile		<50	<50	<50					<50	<50	<50			
Benzene	<5.0	<5.0	<5.0	<5.0	<5.0	<5.0	<5.0	<5.0	<5.0	<5.0	<5.0	<5.0	<5.0	<5.0
Bromobenzene		<5.0	<5.0	<5.0					<5.0	<5.0	<5.0 <5.0			
Bromochloromethane		<5.0	<5.0	<5.0 <5.0					<5.0 <5.0	<5.0 <5.0	<5.0			
Bromodichloromethane Bromoform		<5.0 <5.0	<5.0	<5.0 <5.0					<5.0 <5.0	<5.0	<5.0			
Bronnethane		<10.0	<10.0	<10.0					<10.0	<10.0	<10.0			
n-buty/benzene		<10.0	<10.0	<10.0					<10.0	<10.0	<10.0			
sec-buty@enzene		<10.0	<10.0	<10.0					<10.0	<10.0	<10.0			
fert-butylicenzene		<10.0	<10.0	<10.0					<10.0	<10.0	<10.0			
Carbon disulfide		<10.0	<10.0	<10.0					<10.0	<10.0	<10.0			
Carbon tetrachloride		<5.0	<5.0	<5.0					<5.0	<5.0	<5.0			
Chlorobenzene		<5.0	<5.0	<5.0					<5.0	<5.0	<5.0			
Chlorodibromomethane		<5.0	<5.0	<5.0					<5.0	<5.0	<5.0			
Cirlor ve thane		<10.0	<10.0	<10.0					<10.0	<10.0	<10.0			
Chloroform		<5.0	<5.0	<5.0					<5.0	<5.0	<5.0 <10.0			
Chloromethane		<10.0	<10.0	<10.0					<10.0 <5.0	<10.0 <5.0	<5.0			
o-chlorotokiene		<5.0	<5.0 <5.0	<5.0 <5.0					<5.0 <5.0	<5.0	<5.0			
p-chiorotoluene 1,2-dibromo-3-chioropropane		<5.0	<5.0	<5.0		<b>—</b>			¥3.0	₹9.0	7.0			
1,2-dibromo-3-chioropropane			<del></del>			<b> </b>								<b></b>
Dibromomethane		<5.0	<5.0	<5.0		<b> </b>			<5.0	<5.0	<5.0			
1,2-dichlorobenzene		<5.0		<5.0					<5.0	<5.0	<5.0			
1,3-dichlorobenzene		<5.0	₹5.0	<5.0		·			<5.0	<5.0	<5.0			
p-dichlorobenzene		<5.0	<5.0	<5.0					<5.0	<5.0	<5.0			
Dichlorofluoromethane														
Dichiorodifluoromethane		<10.0		<10.0					<10.0	<10.0	<10.0			
1,2-dichloroethane		<5.0		<5.0					<5.0	<5.0	<5.6			
1,1-dichlorcethane	<5.0			38.9	13	25.2	63.9	58.1	<5.0	<5.0	<5.0	<5.0	<5.0	<5.0
1,1-dichloroethene	<5.0			<5.0	<u> &lt;5.0</u>	<5.0	6.1	8.20	<5.0	<5.0	<5.0	<5.0	<5.0	<5.0
cis-1,2-dichlarcethene	<u> </u>	<5.0		<5.0					<5.0	<5.0	<5.0 <5.0			ļI
trans-1,2-dichloroethene	ļ	<5.0 <10.0	<u> </u>	<5.0 <10.0				<del> </del>	<5.0 <10.0	<5.0 <10.0	<10.0			
trans-1,4-dichloro-2-butene		<10.0	<10.0	<10.0	├	ļ	<b> </b>	ļ	\$10.0	×10.0	10.0	<del> </del>		
di-isopropylether 1,2-dichloropropane		<5.0	<5.0	<5.0	<b></b>	<del> </del>	<del></del>		<5.0	<5.0	<5.0			
1,3-dichloropropane	<del></del>	<5.0		45.0				<del> </del>	<5.0	<5.0			· >	
2,2-dichloropropane		<5.0		<5.0					<5.0	<5.0	<5.0			
1,1-dichloropropene		<5.0		<5.0		<del>                                     </del>			<5.0	<5.0	<5.0			
1,3-dichloropropene		<10.0		<10.0					<10.0	<10.0	<10.0			
cis-1,3-dichloropropene		<5.0	<5.0	<5.0					<5.0	<5.0	<5.0			
trans-1,3-dichloropropene		<5.0							<5.0					
Ethythenzene		<5.0	<5.0	<5.0					<5.0	<5.0	<5.0			
Heptane									<b></b>			<b> </b>		
2-hexanone		<5.0			ļ				<5.0	<del></del>			<u> </u>	
Hexachlorobutadiena	ļ	<10.0			<b> </b>	<b> </b>		ļ	<10.0			<u> </u>	ļ	ļ
Iodomethane	<del> </del>	<10.0			<del> </del>	<del> </del>			<10.0 <10.0					<u>  </u>
jsopropylbenzene p-isopropyliotuene	<del> </del>	<10.0 <10.0	4,		<del> </del>	<del> </del>	<del> </del> -	<del> </del>	<10.0	<del></del>		<del></del>		<del> </del>
Methylene chloride	<del> </del> -	<5.0			<del> </del>	<del> </del>	<del> </del>		<5.0	<del></del>			ļ	
Methyl ethyl kelone	<del> </del>	<10.0				<del> </del>	<del></del>	<del> </del>	<10.0				<del> </del>	<del> </del>
MTBE		1	1	T	<b></b>	1	l	<del> </del>	1	1	T	1	<del> </del>	
Methyl isobutyl ketone	1	1	T	I			<u> </u>					<u> </u>		!
4-methyl-2-pentanona		<5.0	<5.0	<5.0		L			<5.0		·			
Totrahydrofuran		<5.0	<5.0	<5.0				<u> </u>	<5.0	<5.0	<5.0	ļ	<u> </u>	
Naphthalene		ļ	L	<u> </u>	ļ	ļ	ļ	ļ	ļ		ļ	ļ	<b> </b>	
n-propytoenzena	ļ	<10.0				ļ	ļ	ļ	<10.0				<b></b>	
Styrene	ļ	<5.0				ļ	<del> </del>	<del> </del>	<5.0	<del></del>			<del> </del>	ļ
1,1,1,2-tetrachloroethane	<del></del>	<5.0				<del> </del>	<del> </del>	<del> </del>	<5.0			<u> </u>	<del> </del>	<del> </del>
1,1,2,2-tetrachloroe/hane	<del> </del>	<5.0 <5.0				<del> </del>	<del> </del>	<del> </del>	<5.0 <5.0				<del> </del>	<del> </del> -
Tetrachioroethene	<b>45.0</b>					o <5.0	<5.0	<5.0					<5.1	<5.0
Toluene 1,2,3-trichlorobenzane	*5.0	<5.0				7 (3.1	7	1	<5.0				1	1 3.0
11,2,3-trichiorobenzane	<del> </del> -	<5.0				<del> </del>	<del> </del>	<del> </del>	<5.0				<del> </del>	<del> </del>
1,1,1-trichforoethane	<5.0		+ <del></del>			0 5.7	<5.0	<5.0					<6.	<5.0
1,1,2-bichloroethane	<del> </del>	<5.0				† <del></del>	† <u>-</u> -	1	45.0	· · · · · · · · · · · · · · · · · · ·			1 <b>-</b>	Γ
Trichloroethene	İ	₹5.0				1	1	1	₹5.0				1	1
Trichloroflucromethane	1	<10.0				<u> </u>		<u> </u>	<10.0				L	I
1,2,3-trichloropropane	L	<5.0				$\Box$			<5.0	<5.				L
1,2,4-trimothylbenzane		<10.0							<10.0	·				
1,3,5-trimethylbenzene		<10.0					L		<10.0				ļ	
Vinyl acetala		<10.0	<10.0	<10.0	<b> </b>	ļ	<u> </u>	<u> </u>	<10.0	<10.0	<10.0	4	<b></b>	<b></b>
Ethyl acetate		<u></u>	ļ		<del></del>	<b> </b>		ļ	1	ļ	<del> </del>		ļ	<u> </u>
Vinyl chloride		<10.0	<10.0	≰10.0		ļ	ļ	ļ	<10.0	<10.	<10.		ļ	ļ
2-chicroethylvinyl ether	ļ		<del> </del>	<del> </del>	ļ	<del> </del>	<del> </del>	<del> </del>	<del> </del>	<del> </del>	<del> </del>		<del> </del>	<del> </del>
Ethyl methacrylate	ļ	ļ		<10.0		<del> </del>	<del> </del>	<del> </del>	<10.0	<10.	<10.	<del> </del>	<b>+</b> -	<del> </del>
o-xylene	ļ	<10.0 <5.0				<del> </del>	<del> </del>	<del> </del>	<10.0 <5.0				<del> </del>	<del> </del>
rn and p sylene	<del> </del>	<5.0 <5.0				<del> </del>	<del> </del>	<del> </del>	<5.0 <5.0				<del> </del>	<del> </del>
Xylenes, total	L	<u> </u>	< 5.0	L	L	J	I		<p.(< td=""><td>·1</td><td>×1</td><td><u> </u></td><td>1</td><td></td></p.(<>	·1	×1	<u> </u>	1	

Compound	MW8-S	MW9-8	MW9-S	MW9-S	MW9-S	MW9-S	MW9-B	MWS-S	MW1-D	MW1-D	MW1-D	MW1-D	MWI-D	MW1-D
	11/30/00	7/26/99	9/9/99	11/1/99	3/2/00	6/23/00	9/29/00	11/30/00	7/24/98	10/26/98	2/16/99	7/26/99	9/9/99	11/1/99
Acetone		<50	<50	<50					<20	<20		<50	<50	<50
Acroieln		<50	<50	<50					<10	<50		<50 <50	<50 <50	<50 <50
Acrylonitrile		<50	<50	<50					<10	<10 <0.5	<0.5	<5.0	<5.0	<5.0
Benzene	<5.0	<5.0	<5.0	<5.0	<5.0	<5.0	<5.0	<5.0	<0.5 <1.0	<1.0	<1.0	<5.0	<5.0	<5.0
Bromobenzene		<5.0	<5.0 <5.0	<5.0 <5.0					<1.0	<1.0	<1.0	<5.0	<5.0	<5.0
Bromochloromethane		<5.0	<5.0	<5.0 <5.0					<1.0	<1.0	<1.0	<5.0	<5.0	<5,0
Bromodichloromethane		<5.0 <5.0	<5.0	<5.0					<2.0	<2.0	<2.0	<5.0	<5.0	<5.0
Bromoform Bromomethane		<10.0	<10.0	<10.0					<4.0	<4.0	<4.0	<10.0	<10.0	<10.3
n-butyfbenzene		<10.0	<10.0	<10.0					<1.0	<1.0	<1.0	<10.0	<10.0	<10.0
sec-buty/banzene		<10.0	<10.0	<10.0					<1.0	<1.0	<1.0	<10.0	<10.0	<10.0
tert-buty/benzene		<10.0	<10.0	<10.0					<1.0	<1.0	<1.0	<10.0	<10.0	<10.0
Carbon disulfide		<10.0	<10.0	<10.0					<1.0	<1.0		<10.0	<10.0	<10.0
Carbon tetrachloride		<5.0	<5.0	<5.0					<1.0	<0.3	<0.3	<5.0	<5.0	<5.0
Chlorobenzene		<5.0	<5.0	<5.0					<1.0	<1.0	<1.0	<5.0	<5.0	<5.0
Chlorodibromomethane		<5.0	<5.0	<5.0					<1.0	<1.0	<1.0	<5.0	<b>45.0</b>	<5.0
Chloroethane		<10.0	<10.0	<10.0					<4.0	<4.0	<4.0	<10.0	<10.0	<10.0
Chloroform		<5.0	<5.0	<b>&lt;</b> 5.0					<1.0	<1.0	<1.0	<5.0	<5.0	
Chloromethane		<10.0	<10.0	<10.0					<1.0	<1.0	<1.0	<10.0	<10.0	
o-chlorotoluene		<5.0	<5.0	<5.0					<1.0	<1.0	<1.0	<\$.0	<5.0	
p-chiarotoluene		<5.0	<5.0	<5.0					<1.0	<1.0		<5.0	<5.0	<5.0
1,2-dibromo-3-chloropropane									<10	<10				ļ
1,2-dibromosthane									<10	<10			<u></u>	ļ <u>.</u>
Dibromomethane		<5.0	<5.0	<5.0		<b></b>		ļ	<1.0	<1.0			<5.0 <5.0	
1,2-dichlorobenzene		<5.0	<5.0	<5.0		ļ			<1.0	<1.0				
1,3-dichlorobenzene		<5.0	<5.0	<5.0	<u> </u>	ļ	<b> </b>	<u> </u>	<1.0	<1.0			<5.0 <5.0	
p-dichlorobenzene	<u> </u>	<5.0	<5.0	<5.0	<del> </del>	<del> </del>	<del> </del>	<b></b>	<1.0	<1.0		1 5.0		1-3
Dichlom@uoromethane	ļ					<del> </del>	<del> </del>	<b></b>	<10 <3.0	<3.0		<10.0	<10.0	<10.0
Dichlorodifluoromethane	<b></b>	<10.0	<10.0 <5.0	<10.0 <5.0	<b> </b>	<del> </del>		<del> </del>	<1.0	<1.0				4
1,2-dichloroethane	<5.0	<5.0 <5.0			<5.0	<5.0	<5.0	<5.0	<1.0	<5.0				
	<5.0	<5.0	<5.0	<5.0	<5.0			<5.0	42.0					
1,1-dichloroethene cis-1,2-dichloroethene	C3.U	<5.0	<5.0	<u> </u>		1	1	<del></del>	<1.0	<1.5		+		
trans-1,2-dichlorosthene		<5.0	<5.0		<del> </del>	<del> </del>	<del> </del>		<1.0				<del></del>	
trans-1,4-dichloro-2-butene		<10.0	<10.0			<del> </del>	<del> </del>	<del> </del>	<10					
di-isopropylether			1,0.0	1		<del>                                     </del>	<del>                                     </del>		<1.0	<1.0		1	<b></b>	1
1,2-dichloropropane		<5.0	<5.0	<5.0	ļ	<del> </del>	†		<1.0	<1.0		<5.0	<5.0	<5.0
1,3-dichloropropane		<5.0	<5.0				<b> </b>		<1.0	<1.0	<1.0	<5.0	<5.0	<5.
2,2-dichloropropane		<5.0			1		1		<1.0	<1.0	<1.	<5.0	<5.0	<5.
1,1-dichloroproperæ		<5.0		<5.0		1		1	<1.0	<1.0	<1.0	<5.0	<5.0	45.
1,3-dichloropropene		<10.0	<10.0	<10.0		1						<10.0	<10.0	<10.
cis-1,3-dichloropropene		<5.0	<5.0	<5.0					<1.0	<1.	<1.			
trans-1,3-dichloropropene		<5.0							<1.0					
Ethylbenzene		<5.0	<5.0	<5.0			<u> </u>		<1.0		<1.	0 <5.0	<5.0	o <5.
Heptane							<u> </u>	<u> </u>	<1.1	<del></del>	<b></b>		<u> </u>	
2-hexanone		<5.0				ļ	ļ		<10	4		45.0	<del></del>	
Hexachlorobutadiane		<10.0				ļ	ļ		<5.0	<del></del>	<del></del>	<10.0		
lodomethana		<10.0				<b></b>	<del> </del>	ļ	<11		<del></del>	<10.0	4	
Isopropylbenzene		<10.0				<del> </del>		<del> </del>	<1.0				<del></del>	
p-isopropyttoluerie	<b>}</b> -	<10.0				ļ	<del> </del>	<del> </del>	<1.					
Methylene chloride	<del> </del>	<5.0 <10.0				<del> </del>	<del> </del>	<del> </del>	<1		<del></del>	<10.0		
Methyl ethyl ketone MTBE	<del> </del>	<10.0	×10.0	1	<del> </del>	<del> </del>	<del> </del>	<del> </del>	<1.		<del></del>	*10.0	1 -,0.	<del>\ \\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\</del>
Methyl isobutyl xstone	<del> </del>	<del> </del>	<del> </del>	<del> </del>	<del> </del>	<del> </del>	<del> </del>	<del> </del>	<1			<del> </del>	1	1
4-methyl-2-pentarone	<del> </del>	<5.0	<5.0	<5.0	1	<del> </del>	1	1	<del>                                     </del>	<del>                                     </del>	1	<5.0	45.	0 <5
Tetrahydrofuran	<del> </del>	45.0				1	1	1	<5	0 <5	ol .	<5.0		
Naphthalene		†	T		1	T	T	1	<5.				I	1
n-propy/banzene	<del></del>	<10.0	<10.0	<10.0					<1.				<10.	0 <10
Styrene	<u> </u>	<5.0					L		<1.	0 <1.	0 <1	.0 <5.0		
1,1,1,2-tetrachioroethane		<5.0	<5.0			1			<1.					
1,1,2,2-tetrachloroethane		<5.0							<1.					
Tetrachioroethene		<5.0				1			<1.					
Toluena	<5.0					0 <5.	0 15	<5.0						
1,2,3-trichlorobenzene	<u></u>	<5.0				<del> </del>		ļ	< 5.					
1,2,4-trichlorobenzene	<b> </b>	<5.0				<del></del>	<del> </del>	<del> </del>	<5.					
1,1,1-trichioroethane	<5.0					0 <5.	0 <5.	) <u>&lt;5.</u>						
1,1,2-trichloroethane	<b> </b>	<5.0				<del> </del>	<del></del>		<1.					
Trichloroothene	ļ	<5.0				<del> </del>	<del> </del>	<del> </del>	4					
Trichlorofluoromethane	ļ	<10.0				+	<del> </del>	<del> </del>	<1. <1.					
1,2,3-trichloropropane	<del> </del>	<5.0				+	<del> </del>	<del> </del>	1 41					
1,2,4-trimethy/benzene	<del> </del>	<10.0				<del> </del> -	<del> </del>	<del> </del>	- <u>{1</u>	<del></del>				
1,3,5-trimethylbenzene	<del> </del>	<10.0						<del> </del>	- 1			.0 <10. <10.		
Vingi acetate	<del> </del>	<10.0	*10.0	× 10.1	<del>' </del> -	<del> </del>	+	<del> </del> -	<del>                                     </del>	<del>~~~</del>	<del>'1</del>	1 10.	<del>* </del>	<del></del>
Ethyl acetate	<del> </del>		<10.0	<10.	,	+	<del> </del>	<del> </del>	- 41	.0 -1	.0 <1	.0 <10	0 <10	.0 <10
Vinyl chloride		<10.0	*10.0	<10.	1	+	+	<del> </del>	<del> *</del>	<del>" "</del>	<del>~</del>	- C10	<del></del>	* * * * * * * * * * * * * * * * * * * *
2-chloroothylvinyl ether	<del> </del>		<del> </del>	<del> </del>	1	+	+	<del> </del>	+	+	+	+	<del>- </del>	<del>- </del>
Ethyl methacrylata	<del> </del>	<10.0	<10.0	×10.		+	<del></del>	+	+	+	+	<10	.0 <10	.0 <1
o-xylene	ļ	<10.0 <5.0				<del>- </del>	<del></del>	<del> </del>	+	<del> </del>	+	<5	~~~~~	
m and p xylene				-, -,,,	- 1		1							

					16155	1949 7	10100	inin n 1	inin n	APID D	ann n I	MW2-D	MW2-D	MW3-D
Cumpourd	MW1-D	MW1-D	MW1-D	MW1-D	MWZ-D	MW2-D	MW2-D	MW2-D	MW2-D	MW2-D	MW2-D		1	7/24/98
	3/2/00	6/23/00	9/29/00	11/30/00	7/24/98	10/26/98	7/26/99	9/9/99	11/1/99	3/2/00	6/23/00	9/29/00	11/30/00	1/2430
Acetone							<50 c50	<50 <50	<50 <50					
Acrolein Acrylonitrile							<50 <50	<50	<50 <50					
Benzene	<5.0	<5.0	<5.0	<5.0	<0.5	<0.5	<5.0	<5.0	45.0	<5.0	<5.0	<5.0	<5.0	40.5
Bromobenzene		10.0				-0,0	<5.0	<5.0	<5.0					
Bromochlorometrane				~			<5.0	<5.0	<5.0					
Bromodichloromethane							<5.0	<5.0	<5.0					
Bromoform							<5.0	<5.0	<5.0					
Bromomethane							<10.0	<10.0	<10.0					
n-butytbenzene							<10.0	<10.0	<10.0					
sec-butyfbenzene							<10.0	<10.0	<10.0					
tert-butyfbenzene							<10.0	<10.0	<10.0					
Carbon disulfide							<10.0	<10.0	<10.0					
Carbon tetrachloride							<5.0	<5.0	<5.0					
Chlorobeitzene		I					<5.0	<5.0	<5.0					
Chlorodikromomelitane	ļ	<u> </u>					<5.0	<5.0	<5.0					
Chloroethane							<10.0	<10.0	<10.C <5.0					
Chloroform	ļ	<u> </u>		<del></del>			<5.0	<5.0 <10.0	<10.0					*********
Chloromethana		<del> </del>					<10.0	<5.0	<5.0					
o-chiorololuene p chiorololuene	<del>                                     </del>	<del> </del>	<b> </b>				<5.0 <5.0	<5.0	<b>45.0</b>				<del></del>	
1,2-dibromo-3-chloropropane	<del> </del>	<b></b>	<b></b>	<del></del>	<b></b>				~5,0			·		
1,2-dibromoethane	<del> </del>	<del> </del>	<del> </del>	<del> </del>	<b></b>		<b></b>						<del></del>	
Dibromomethane	<del>                                     </del>	<del>                                     </del>	l	····	<del> </del>	<del></del>	<5.0	<5.0	<5.0					
1,2-áchiorobenzene	<del>                                     </del>	<u> </u>	<b></b>		<del> </del>	<b></b>	<5.0		<5.0					
1,3-dichlorobenzene	†						<5.0		<5.0					
p-dichlorobenzene					<u> </u>		<5.0	<5.0	<5.0					
Dichlorofluoromethane														
Dichlorodifluoromethane	1	L					<10.0	<10.0	<10.0					
1,2-dichloroethane		L					<5.0	<5.0	<5.0	<b></b> i				
1,1-dichloroethane	<5.0			<5.0		ļ	<5.0	<5,0	<5.0		<5.0	<5.0	<5.0	
1,1-dichloroethene	<5.0	<5.0	<5.0	<5.0	<2.0		ļ	<5.0	<5.0	<5.0	<5.0	<5.0	<5.0	<2.0
cis-1,2-dichioroethene	<u> </u>	<del> </del>		<b></b>			ļ	<5.0	<5.0					
trans-1,2-dichloroethene	<del> </del>	<del> </del>	ļ		ļ		ļ	<5.0	<5.0 <10.0					
trans-1,4-dichloro-2-butene di-Isopropytether	<del> </del>	ļ		<del></del> -	<del> </del>	<del></del>	<del> </del>	<10.0	¥10,0					
1,2-dichloropropane	<del> </del>	<del> </del>		<del> </del>			<del> </del>	<5.0	<5.0					
1,3-dichloropropane	┼──	<del> </del>	<del></del>	<del></del>	<del></del>		<del> </del>	<5.0	<5.0					
2,2-dichloropropane	<del> </del>	<del>                                     </del>	<del>                                     </del>					<5.0	<5.0	<del> </del>	<del></del>			
1,1-dichloropropene	<del> </del>	<b></b>		<del> </del>	<del> </del>		<u> </u>	<5.0	<5.0					
1,3-dichloropropene	<del>                                     </del>	<b> </b>	<b>—</b>					<10.0	<10.0					
cla-1,3-dichloropropena								<5.0	<5.0					
trans-1,3-dichloropropone	Ĺ							<5.0	<5.0					
Ethythenzene	<u> </u>		<u> </u>	ļ		<u> </u>		<5.0	<5.0					
Heptane	ļ	ļ		<u> </u>	<del> </del>		ļ		ļ <u>.</u>			ļi		
2-hexanone		<del> </del>	<del></del>	<del> </del>	<del> </del>		<u> </u>	<5.0	<5.0					<b></b>
Hexachixrobutatione	ļ	<del> </del>		ļ	<del> </del>	<del> </del>	<del> </del>	<10.0 <10.0	<10.0 <10.0					
lodomethane isopropytbenzene	<del> </del>	<del> </del>		<del> </del>	<del> </del>	<del> </del>	<del> </del>	<10.0	<10.0					
p-Isopropyltokiene	<del> </del> -	<del> </del>	<del></del>	<del></del>	<del> </del>	<del> </del>	<del> </del>	<10.0						
Methylene chicride	<del> </del>	┼──	<del> </del>		<del> </del>	<del>                                     </del>	<del> </del>	<5.0			<b>——</b>			
Mothyl ethyl ketone	<del> </del>	<b>†</b>	1	1	1		1	<10.0	<10.0		<b></b>	l		
мтве	1													
MeGryl isobutyl ketone														
4-methyl-2-pentanone	1						L	<5.0						
Tetrahydrofuran	<b></b>	<b> </b>	<b> </b>	ļ		<u> </u>	<b>!</b>	<5.0	<5.0			L	ļ	
Naphthalene	<del> </del> -	<del> </del>	<b> </b>	<b></b>	<b> </b>	<b></b>	<b> </b>	<del> </del>			ļ	ļ	ļ	
n-propylbenzene	ļ	<del> </del>	<del> </del>		<del> </del>	<del> </del>	<del> </del>	<10.0				<b> </b>	<b>_</b>	
Styrene	<del> </del>	<del> </del>	<5.0 <5.0			<del> </del>	<del> </del>							
1,1,1,2-letrachloroethane	<del> </del>	<del> </del>	<5.0 <5.0			<del> </del>	<del> </del>	<del> </del>	<del> </del>					
1,1,2,2-tetrachloroethane Tetrachloroethene	<del> </del>	<del> </del>	<5.0			<del> </del>	<del> </del>	<del></del>	I					
Toluene	<5.0	<5.0	<5.0	<5.0	<1.0	<1.0	<del> </del>	<5.0			<6.0	<5.0	<5.0	<1.0
1,2,3-trichlorobenzene	1 .0.1	1	1	<del> </del>	1		<del> </del>	<5.0			<del>}}</del>	1		`1.0
1,2,4-tilchlorobenzene	<del>                                     </del>	<del> </del>	<del> </del>	<del>                                     </del>	<del>                                     </del>	<del> </del>	<del> </del>	₹5.0			<del> </del>	<del></del>	<del> </del> -	<del> </del>
1,1,1-trichlorcethane	<5.0	<5.0	<5.0	<5.0	<1.0	<1.0	1	<5.0			<5.0	<5.0	<5.0	<1.0
1,1,2-trichioroethane	† <del></del>	1	<u> </u>	1	1	1	1.	<5.0			1	<u> </u>	T	
Trichloroethene	1	<u> </u>	1	1		<1.0		<5.0			1		<u> </u>	<1.0
Trichlorofluoromethane								<10.0			L		L	
1,2,3-trichloropropane								<5.0						
1,2,4-trimethy/benzene								<10.0						
1,3,5-trimethy/benzene		1						<10.0						
Vinyl ecetate						ļ	<u> </u>	<10.0	<10.0					1
Ethyl acetate		ļ	ļ	<u> </u>	ļ	<b> </b>	<u> </u>	ļ		ļ	<u> </u>	ļ		
Vinyl chloride	ļ	<b></b>	ļ	ļ	ļ	<b>↓</b>	<b></b>	<10.0	<10.0	1	ļ	ļ		ļ
2-chloroethylvinyi ether	ļ	<b> </b>	ļ	<del></del>	ļ	<del> </del>	<del> </del>	ļ	<b>ļ</b>	<b>ļ</b>	<del> </del>	<b></b>	<del> </del>	ļ
Ethyl methacrylate	<del> </del>	<del> </del>		ļ	<del> </del>	<del> </del>	<del> </del> -	<del> </del>	·		ļ	<del> </del>	<del> </del>	<del></del>
o-xylene		<b> </b>	<del> </del>	ļ	<del> </del>	<b></b>	<del> </del>	<10.0 <5.0			<del> </del>	<del> </del> -	<del> </del>	ļ
m and p xylene	<del> </del>	<del> </del>		<del> </del>	<del> </del>	<del> </del>	<del> </del>	<5.0 <5.0			<del> </del> -	<del> </del>	<del> </del>	<del> </del>
Xylenes, total	L	<u></u>	L	l:	<del></del>	<del></del> ,	J			1		1	<u></u>	<u> </u>

Compound	MW3-D	MW3-D	MW3-D	MW3-D	MW3-D	MW3-D	MW3-D	MW3-D	MW4-D	MW4-D	MW4-D 7/26/99	MW4-D 9/9/99	NW4-D 11/1/99	MW4-D 3/2/00
	10/26/98	7/26/99	9/9/99	11/1/99	3/3/00	6/23/00	9/29/00	11/30/00	7/24/98	10/26/98		<50	<50	3/2/0
Acetone		<50	<50	<50					<50 <50		<50 <50	<50	- 50	
Acrolein		<50 <50	<50 <50	<50 <50					<b>450</b>		-\$0 -\$0	<b>&lt;50</b>	<50	
Aczylonitriie Benzene	€0.5	<5.0	<5.0	<5.0	<5.0	<5.0	<5.0	<5.0	<b>45.0</b>	<0.5	<5.0	<5.0	<5.0	<5.0
Bromobenzene		<5.0	<5.0	<5.0	V.0	3.0	-0.0	-0.0	₹5.0		<5.0	<5.0	<5.0	
Bromochloromethane		<5.0	<5.0	<5.0					<5.0		<5.0	<5.0	<5.0	
Gromodichloromethane		<5.0	<5.0	₹5.0					<5.0		<5.0	<5.0	<5.0	
Bromoform		<5.0	<5.0	<5.0					<5.0		<5.0	<5.0	<5.0	
Bromomethane		<10.0	<10.0	<10.0					<10.6		<10.0	<10.0	<10.0	
n-butyfbenzene		<10.0	<10.0	<10.0					<10.0		<10.0	<10,0	<10.0	
sec-buty/benzerie		<10.0	<10.0	<10.0					<10.0		<10.0	<10.0	<10.0	
tert-buty/benzene		<10.0	<10.0	<10.0					<10.0		<10.0	<10.0	<10.0	
Carbon disulfide		<10.0	<10.0	<10.0					<10.0		<10.0	<10.0	<10.0	
Carbon tetrachlorido		<5.0	<5.0	<5.0					<5.0		<5.0	<5.0	<5.0	
Chiorobenzene		<5.0	<5.0	<5.0					<5.0		<5.0	<5.0	<5.0	
Chloro@bromomethane		<5.0	<5.0	<5.0					<5.0		<5.0	<5.0	<5.0	
Chloroethane		<10.0	<10.0	<10.0					<10.0		<10.0	<10.0	<10.0	
Chloroform		<5.0	<5.0	<5.0					<5.0		<5.0	<5.0	<5.0	
Chloromethane		<10.0	<10.0	<10.0					<10.0		<10.0	<10.0	<10.0	
o-chicrotoluene		<5.0	<5.0	<5.0		<b></b> _		ļ	<5.0		<5.0	<5.0	<5.0	
p-chiorotoluene		<5.0	<5.0	<5.0		ļ	<u> </u>		<5.0		<5.0	<5.0	<5.0	
1,2-dibromo-3-chloropropane				<u> </u>		<del> </del>		ļ	<b> </b>			<b>  </b>		
1,2-dibromoethane	ļ		ļ <u>-</u> -	<u> </u>	ļ	<del> </del>	<b></b>	<del> </del>				<5.0	<5.0	
Dibromomethane		<5.0	<5.0	<5.0	ļ	<del> </del>	<del> </del>	<del></del>	<5.0		<5.0		<5.0 <5.0	
1,2-dichlorobenzene		<5.0	<5.0 <5.0	<5.0 <5.0	<del> </del>		<del> </del>	<del> </del>	<5.0 <5.0		<5.0	<5.0 <5.0	45.0	
1,3-dichlorobenzene p-dichlorobenzene	<b> </b>	<5.0 <5.0	<5.0 <5.0	<5.0 <5.0	ļ	<del> </del>	<del> </del>	<del></del>	<5.0 <5.0		<5.0 <5.0	<5.0 <5.0	<5.0	
<del></del>		9.0	45.0	(5,0				<del> </del>	19.0			1000	72.0	
Dichlorofluoromethane Dichlorodifluoromethane		<10.0	<10.0	<10.0	<del> </del>	<del> </del>	<del> </del>	<del> </del>	<10.0		<10.0	<10.0	<10.0	
1,2-dichloroethane	<del></del>	<3E	<5.0	<5.0	<del> </del>	<del> </del>	<del>                                     </del>	<del>                                     </del>	<5.0		<5.0	45.0	<50	<del></del>
1,1-dichloroethane	<1.0		<5.0	<5.0		<5.0	<5.0	<5.0	1	<1.0	<5.0		<5.0	<5.
1,1-dichloroethene		<5.0	<5.0	<5.0		<del></del>		<del></del>	<2.0		<5.0		<5.0	<5.
cis-1.2-dichloroethane	<del> </del>	<5.0		<5.0					<u> </u>		<5.0		<5.0	
i ans-1,2-dichloroethene		<5.0	<5.0	<5.0		†	<del>                                     </del>	<u> </u>	<del> </del>		<5.0		<5.0	
trans-1,4-Gichloro-2-butene		<10.0	<10.0	<10.0		1					<10.0	<10.0	<10.0	
di-Isopropylethar		1			<b>1</b>				<b> </b>					
1,2-dichloropropane		<5.0	<5.0	<5.0		1			· · · · · · · · · · · · · · · · · · ·		<5.0	<5.0	<5.0	
1,3-dichloropropane		<5.0	<5.0	<5.0							<5.0	<5.0	<5.0	
2,2-dichioropropane		<5.0	<5.0	<5.0							<5.0	<5.0	<5.0	
1,1-dichloropropene		<5.0	<5.0	<5.0	· · · · ·				Ι.		<5.0	<5.0	<5,0	
1,3-dichloropropene		<10.0	<10.0	<10.0							<10.0	<10.0	<10.0	
cis-1,3-dichloropropene		<5.0	<5.0	<5.0							<5.0		<5.0	
trans-1,3-dichloropropene		<5.0	<5.0	<5.0							<5.0		<5.0	
Ethylbenzerie		<5.0	<5.0	<5.0			ļ		<u> </u>		<5.0	<5.0	<50	
Heptane	<u> </u>		<b></b>		<b> </b>	ļ	<u> </u>	ļ	ļ	ļ		ļ		
2-hexanone		<5.0	<5.0			ļ	<b></b>	<u> </u>	ļ	ļ	<5.0		<5.0	
Hexachlorobutadiene		<10.0	<10.0	<10.0							<10.0	<del></del>	<10.0	ļ <u>.                                 </u>
lodomethane	ļ	<10.0	<10.0 <10.0	<10.0 <10.0		<del> </del>	<del> </del>	<del> </del> -	<del> </del>	<b></b> -	<19.0		<10.0 <10.0	ļ
Isopropylbenzene	<del> </del>	<10.0 <10.0	4	<10.0		<del> </del>	<del> </del>	<del> </del>	<del>}</del> -	<del> </del>	<10.0 <10.0		<10.0	<del> </del>
p-isopropytoluene Methylene chlorida		<5.0	<del></del>	<5.0		<del> </del>	<del> </del>	<del> </del>	<del> </del>		<5.0		<5.0	
Methyl ethyl ketone		<10.0		<10.0		<del> </del>	<del> </del>	<del></del>	<del> </del>	<del> </del> -	<10.0		<10.0	
MTBE	<del> </del>	1 - 2,10.0	1	1 -10.0	<del>                                     </del>	1	<b>-</b>	<del> </del>	1	<del>                                     </del>	1	1	<del>-</del>	<del>                                     </del>
Methyl isobutyl ketone	<del> </del>	1	1		<del> </del>	1	1	<del>                                     </del>	<del> </del>	t	1	1	<del> </del>	<del>                                     </del>
4-methyl-2-pentanone		<5.0	<5.0	<5.0		1	1.	1	1	1	<5.0	<5.0	<5.0	1
Tetrahydrofuran	I	<5.0		<5.0			L		I		¢5.0		<5.0	
Naphthalene									1					
n-propy@enzene		<10.0				1					<10.0			
Styrene		<5.0				1					<5.0			
1,1,1,2-tetrachioroethane		<5.0				<b></b>	ļ	<b></b>	<b> </b>	ļ	<5.0		<5.0	
1,1,2,2-tetrachloroethane		<5.0				<u> </u>	<b> </b>	<b> </b>	<b> </b>	ļ	<5.0			
Tetrachioroathena	ļ	<5.0				<del> </del>	<del></del>	ļ		ļ	<5.0			
Toluene	<1.0					0 <5	0 <5.0	<5.0	<1.0	<1.0				
1,2,3-trichlorobenzene		<5.0				<del> </del>	<del> </del>	<del> </del>	<del> </del>	ļ	<5.0			
1,2,4-trichlorobenzene	<1.0	<5.0 <5.0				o <5.	0 <5.0	<5.0	<1.0	<1.0	<5.0			
1,1,1-trichloroethane	\$1.0	<5.0				<del>`````</del>	7	1	<del>' </del>	<del>                                     </del>	×5.1			
<del></del>	<1.0					<del> </del>	<del> </del>	+	<1.0	×1.0				
Trichloroethene Trichlorofluoromethane	<del> </del>	<10.0				<del> </del> -	1	<del> </del>	+	<del> </del>	<10.			
1,2,3-trichloropropane	†	<5.0				<del> </del>	<del> </del>	<del> </del>	1	<del> </del>	<5.			
1,2,4-bimethybenzene	<del>                                     </del>	<10.0				1	<del> </del>	<del> </del>	┼──	<del> </del>	<10.			
1,3,5-timethy/benzene	<del> </del>	<10.0				<del> </del>	<del> </del>	<del> </del>	1	<del> </del>	<10			
Virtyl acetate	<del> </del>	<10.0	<del></del>			<del> </del>	1	<del> </del>	<del> </del>	1	<10		4	
Ethyl acetata	<del>                                     </del>	1	l	† <del></del>	†	<del> </del>	<del> </del>	<del> </del>	1	<del> </del>	1	1	1	1
Vinyl chkride	<del> </del>	<10.0	<10.0	<b>&lt;10.0</b>	1	1	<del>                                     </del>	<del> </del>	†	<del> </del>	<10	0 410.0	<10.0	<del> </del>
2-chloroethylvinyl ether	<del> </del>	1	1	T	1	<del>1</del>	<del>                                     </del>	<del> </del>	<del> </del>	<del> </del>	1	1	<del> </del>	
Ethyl methacrylate	<del> </del>	<del> </del>	1	<del> </del>	<del>                                     </del>	1	<del> </del>	<del>                                     </del>	1	1	1	<u> </u>	<del> </del>	<del> </del>
O-XYISHO	<del> </del>	<10.0	<10.0	<10.0	1	1	1	<del> </del>	1	1	<10.	0 <10.0	<10.0	<del>,</del>
m and p xylene	<del> </del>	<5.0	<del></del>			<del> </del>	†	1	1	<del> </del>	<5.			
			<5.0				<del> </del>	+		****	<5.			<del></del>

March   Marc	Compound	MW4-D	MW4-D	MW4-D	MW5-D	KW5-D	MWS-D	MW5-D	MW5-D	MW5-D	MW5-D	MWE-D	MW6-D	MW6-D	kfW8-D
Legister	***************************************			ĺ											
Section	Acetone		- 373.00	1,,,,,,,,				3200	41410	312000	11100175				
Incompany (40 40 40 40 40 40 40 40 40 40 40 40 40 4	Acrolein														
Marcelaremen	Acrytonitrile														
Secondaries	Benzene	<5.0	<5.0	<5.0				<5.0	<5.0	<5.0	<5.0				<5.0
Street Control Process			i												
150   150															
100   1700   1	Bromoform														
100   100	Bromomethane														
enterhybenome	n-but/iberizene				<10.0		<10.0					<10.0			
Carbon bands	sec-butythenzene														
Carlon terralskis  Chronistomore	<u> </u>														
Coloradormembare	The second secon		<u> </u>							<u> </u>					——
1-000   1-00		l													
Colorents	Chlorodibromomethane		<b>—</b> —												
Characterium	Chloroethane				<10.0	<10.0	<10.0					<10.0	<10.0	<10.0	
1,500   1,50	Chloroform														
Commonshame			ļ												
1,280mon-bare			<u> </u>							<b></b>	<del></del>				
3.486monthurs		<del>                                     </del>			<5.0	<5.0	<5.0					<5.0	<5.0	<5.0	
Description		<del> </del>	<del> </del>	<b></b>				<del> </del>				ļ			
1.24-discretareame	Dibromomethane				<5.0	<5.0	<5.0					<5.0	<5.0	<5.0	
\$60   \$60	1,2-dichlorobenzene	L				<5.0	<5.0							<5.0	
Chichostonoreshape	1,3-dichlorobenzene														
Chitocologramma		ļ	<u> </u>		<5.0	<5.0	<b>&lt;</b> 5.0					<5.0	<5.0	<5.0	
12-601 tonorwine		<del> </del> -	<del> </del>	<b></b>			740.0			<u> </u>					
1.1.64/bronomer		<del> </del>		<b></b>											
1.5de/broorbane	1,1-dichloroethane	<5.0	<5.0	<5.0				<5.0	<5.0	<5.0	<5.0				<5.
Image   1,50	1,1-dichloroethene														
Table   1450hr   2040hr   100   10	cts-1,2-dichloroethene				<5.0	<5.0						<5.0	<5.0	<5.0	
dispropryshed	trans-1,2-dichloroothene														
13-delingrogroups			<b></b> _		<10.0	<10.0	<10.0					<10.0	<10.0	<10.0	
13-def-harpopropium   \$40		<del> </del>	<del> </del>				-E 0								
23 del Arrogrophies															
1.1.doi:1000000000000000000000000000000000000		t													
displayments	1,1-dichloropropene														
	1,3-dichloropropene				<10.0							<10.0	<10.0	<10.0	
Ethystorians   150   150   150   150   150   150	cis-1,3-dichloropropene														
Page		ļ	ļ												
2-bas anone   4.50					<5.0	\$.0	*5.0	<del> </del> -				<5.0	<5.0	<5.0	
Hastachiootadadene	2-hexanone	<del>                                     </del>			<50	€.0	<5.0					<5.0	<50	<50	
Inspiration	Hexachlorobutadiene		·												
Pipping	Indomethane				<10.0	<10.0	<10.0					<10.0	<10.0	<10.0	
Methylene chizoide	lscpropy/benzene	<u></u>													
Methy statistics			ļ												
Metry  Second		<del> </del>	<del></del>					ļ		<b></b>					<u></u>
Methylisobory  iselone	MENYI ENIYI KOTONE		<del> </del>	<b></b>	-10.0		10.0	t			<del></del>	×10.0	<10.0	<10.0	
Company   Comp	Methyl isobutyl ketone		l			l		· · · · · · · · · · · · · · · · · · ·			<u> </u>		<u> </u>		
Naphthatene	4-metryl-2-pentanone				<5.0	<5.0	<5.0					<5.0	<5.0	<5.0	[
Stylene	Tetrahydrofuran				<5.0	<5.0	<5.0					<5.0	<5.0	<5.0	
Syrene		<b> </b>	ļ					ļ		ļ			<b></b>		
1.1.1.2-tetrachiorosthane			<del> </del>							ļ	ļ				}
1,1,2.24stachbroethane	1,1,1,2-letrachloroethane									<del> </del>					<b></b>
Tetachtoroethens	1,1,2,2-tetrachloroethane										<b></b>				
Tolume	Tetrachloroethena				<5.0	<5.0	<50								/i
1.2.4 techtorobenzane	Toluene	<5.0	<5.0	<5.0					<50	<50	<5.0	<5.0	<5.0	<5.0	<5.
1,1,1-sk/hicroethane										ļ					
1,1,2-tichiprocethane															
Firchizorothans		*5.0	*5.0	<b>*3.0</b>					C3.0	*5.0	*5.0				<u> </u>
Frichtrofluoronethane	Trichicroethene	<b></b>		·					<del> </del>	<del> </del>	<del> </del>				
1,2,3 tichkropropane	Trichlorofluoromethane									<del> </del>	·				<del> </del>
1.2.4 timethy/benzane	1,2,3-trichloropropane				<5.0	<5.0	<5.0			L	<u> </u>				
First script	1,2,4-trimethy/benzene											<10.0	<10.0	<10.0	
Etyl scalula	1,3,5-trimethyfbenzene										<u> </u>		•		
/ary/ chroride	Vriyl acetate	<u>  </u>			<10.0	<10.0	<10.0				ļ	<10.0	<10.0	<10.0	
Criticosthylvinyl ether									ļ	ļ	<b> </b>	<del></del>		ļ	
Etyl methacytate					<10.0	<u></u>	<u> </u>	·		ļ	<del> </del>	<10.0	<10.0	<10.0	
>-ylene								<b></b>		<del> </del>	<del> </del>	ł	<del> </del>	<b> </b>	
and p sylene <50 <50 <50 <50 <50 <50 <50 <50	o-zylene				<10.0	<10.0	<10.0		<del></del>		<del> </del>	<100	<10.0	<100	i
	m and p xylene										L				
	Xylones, total				<5.0	′ <50	<50								

Compound	MW6-D	MW8-D	MW6-D	MW7-D	MW7-D	MW7-0	MW7-D	MW7-D	MW7-D	MW7-D	MW8-D	HW8-D	MW8-D	G-3WM
Compound	6/23/00	9/29/00	11/39/00	1/26/99	9/9/99	11/1/99	3/2/00	6/23/00	9/29/00	11/30/00	7/26/99	9/9/99	11/1/99	3/2/00
Acelone			- 11/100	<50	<50	<50					<50	<50	<50	
Acrolein				<50	<50	<50					<50	<50	<50	
Acrylonitrile				<50	<50	<50					<50	<50	<50	
Benzene	<5.0	<5.0	<5.0	<5.0	<5.0	<5.0	<5.0	<5.0	<5.0	<5.0	<5.0	<5.0 <5.0	<5.0 <5.0	<b>&lt;</b> \$.0
Bromobenzene				<5.0	<5.0	<5.0			<b> </b>		<5.0 <5.0	<5.0 <5.0	<5.0 <5.0	
Bromochloromethane Bromodichloromethane		<b></b>		<5.0 <5.0	<5.0	<5.0 <5.0					<5.0	<5.0	<5.0	
Bromoform				<5.0	<5.0	<5.0					<b>&lt;</b> 5.0	<5.0	<5.0	
Bromomethane				<10.0	<10.0	<10.0					<10.0	<10.0	<10.0	
n-buty <b>ibenzene</b>				<10.0	<10.0	<10.0					<10.0	<10.0	<10.0	
sec-bulylbenzene				<10.0	<10.0	<10.0					<10.0	<10.0	<10.0	
tert-butyfbenzene		ļi		<10.0	<10.0	<10.0					<10.0	<10.0	<10.0	
Carbon disulfide		ļl		<10.0	<10.0	<10.0					<10.0 <5.0	<10.0 <5.0	<10.0 <5.0	
Carbon tetrachloride				<5.0 <5.0	<5.0 <5.0	<5.0 <5.0		<b></b>			<5.0	<5.0	<5.0	
Chlorodatromomethane				<5.0	45.0	<5.0					<5.0	<5.0	<5.0	
Chloroethane				<10.0	<10.0	<10.0					<10.0	<10.0	<10.0	
Chloroform				<5.0	<5.0	<5.0					<5.0	<5.0	<5.0	
Chloromethane				<10.0	<10.0	<10.0					<10.0	<10.0	<10.0	
o-chlorojoluene				<5.0	<5.0						<5.0	<5.0	<5.0	
p-chlorololuene				<5.0	<5.0	<5.0					<5.0	<5.0	<5.0	
1,2-Sibrome-3-chloropropane		ļ	<del></del>		ļ	ļ		<b></b>	<b> </b>		<b> </b>		ļ	
1,2-dibromoethane Dibromomethane		<b></b> -		<5.0	<b>45.0</b>	<5.0		<del> </del>	<b> </b>		<5.0	<5.0	<5.0	<b></b>
1,2-dichlorobenzene		<b></b>		-5.0	<5.0	<5.0		<del>                                     </del>			<5.0	<5.0	<5.0	<del> </del>
1,3-dichlorobenzene		<del></del>		<5.0	<5.0	<5.0	<del> </del>				<5.0	<5.0	<5.0	
p-dichlorobenzens				<5.0	<5.0	<5.0					. <5.0	<5.0	<5.0	
Dichlorofluorometiane								ļ	ļ					
Dichlorodifluoromethane				<10.0		<10.0		ļ	ļ	<u> </u>	<10.0	<10.0	<10.0	ļ
1,2-dichloroethane	ļ			<5.0	<5.0	<5.0		J	<5.0	-50	<5.0	<5.0 <5.0	<5.0 <5.0	
1,1-dichloroefnane	<5.0 <5.0		<5.0 <5.0	<5.0 <5.0	<5.0		<5.0		45.0 45.0	<5.0 <5.0	<5.0 <5.0	<5.0	<5.0	<5. <5.
1,1-dichloroethene cis-1,2-dichloroethene	30.0	3.0	<b>\\$3.0</b>	<5.0			3.0				<5.0		<5.0	<del></del>
trans-1,2-dichloroethene	l	<del>                                     </del>		<5.0				<del> </del>	<del> </del>		<5.0	<5.0	<5.0	
trans-1,4-dichloro-2-butane				<10.0	<10.0						<10.0	<10.0	<.0.0	
di-isopropylether														
1,2-dichloropropane			·	<5.0							<5.0	<5.0	<5.0	
1,3-dichloropropane				<5.0				<u> </u>	ļ		<5.0	<5.0	<5.0	
2,2-dichioropropane	ļ	<b></b>	ļ	<5.0				<b>├</b>	<b>├</b> ──		<5.0	<5.0	<5.0	
1,1-dichioropropene	<del> </del>		<b></b>	<5.0 <10.0			<del> </del>	<del> </del>	<del> </del>		<5.0 <10.0	<5.0 <10.0	<5.0 <10.0	
1,3-dichloropropene cis-1,3-dichloropropene	<del> </del>	<del></del>	<del> </del>	<5.0			<del> </del>	<del> </del>			<5.0	<5.0	<5.0	_
trans-1,3-dichloropropene	<del> </del>	<del> </del>	<del> </del>	₹5.0			<del> </del>	<del> </del>	<del></del>		<5.0	<5.0	<5.0	
Ethylbenzene		<u> </u>		<5.0			<del> </del>	1	<del></del> -		<5.0	<5.0	<5.0	
Heptane														
2-hexanone				<5.0							< 5.0	<5.0	<5.0	
Hexachiorobutadiene		<del> </del>	<b>ļ</b>	<10.0				<del> </del>	<del></del>	ļ	<10.0		<10.0	
lodomethane		<del> </del>	<del> </del>	<10.0 <10.0		4		<del> </del>	ļ	<del> </del>	<10.0 <10.0		<10.0 <10.0	1
Isopropythenzene p-tsopropyttoluene		<del> </del>	<del> </del>	<10.0			<del>}</del> -	<del> </del>	<del> </del>	<del> </del>	<10.0			+
Methylene disoride		<del></del>		<5.0		<del></del>	<del> </del>	<del>                                     </del>			<b>&lt;</b> 5.0		<5.0	
Methyl etryl kelone		1		<10.0				1	1		<10.0	<10.0	<10.0	
MTBE														
Mathyl Isobutyl ketone		ļ	L					ļ	<u> </u>		<u> </u>			
4-methyl-2-pentanone	<del> </del> -	<del> </del>	<del>                                     </del>	<5.0				<del> </del>	<del> </del>	<del> </del>	<5.0			
Tetrahydrofuran Naphthalene	1	<del> </del>	<del> </del>	<5.0	×3.0	<5.0	<del> </del>	<del> </del>	<del> </del>	<del> </del>	<5.0	<5.0	<5.0	<del> </del>
n-propyibenzene	1	<del>                                     </del>	<del> </del> -	<10.0	<10.0	<10.0		1	<del> </del>		<10.0	<10.0	<10.0	<del> </del>
Styrene	L			<5.0	<5.0	<5.0					<b>45.0</b>			
1,1,1,2-tetrachloroethane				<5.0							<5.0			
1,1.2,2-tetrachloroethane		<u> </u>	ļ	<5.0				<del></del>	<b></b>	<b></b>	<5.0			
Tetrachloroetiene	<u> </u>	<del></del>	ļ	<5.0					<del></del>	<del> </del>	<5.0			<del></del>
Toluene 1.2.3-trichlorobenzene	<5.0	<5.0	<5.0	<5.0 <5.0				0 <5.0	o <5.0	<5.0	<5.0 <5.0			
1,2,4-trichkrobenzene		<del> </del>	<del> </del>	<5.0				<del> </del>	<del> </del>	<del> </del>	<5.0			
1,1,1-trichloroefiane	<5.0	<5.0	<5.0					d <5.0	0 <5.0	<5.0				
1,1,2-trichloroethane			L	<5.0	<5.0	<5.0			L	I	<5.0			
Trichicroethene				<5.0	<5.0						<5.0	<5.0	<5.0	
Trichkrofluoromethane	ļ			<10.0					<u> </u>	ļ	<10.0			
1,2,3-trichloropropane	ļ	<del></del>	<b></b>	<5.0				<b></b>	<b>↓</b>	<del> </del>	<5.0			
1,2,4-trimethy/benzene	<del> </del>	<del></del>	<del> </del>	<10.0				<del> </del>	<del> </del>	<b></b>	<10.0			
1,3,5-trimethy/benzene	<del> </del>	<del> </del>	<del> </del>	<10.0 <10.0				<del> </del>	<del> </del>	<del> </del>	<10.0			
Virryl acetate Ethyl acetate	<del> </del>	<del> </del>	<del> </del>	×10.0	1 10.0	*10.0	<del> </del>	+	+		<10.0	*10.0	×10.0	<del>' </del>
Vinyl chioride	<del> </del>	†	<del> </del>	<10.0	<10.0	<10.0	1	†	<del> </del>	+	<10.0	<10.0	<10.0	d
2-chloroethylvinyl ether		1	<u> </u>	1	1	1	<del> </del>	1	1	t	1	1	1	1
Ethyl methacrylate	1	1	I		<u> </u>	1	<u> </u>		L	<u> </u>			1	1
	1	T	1	<10.0	<10.0	<10.0					<10.0	<10.0	<10.0	0
o-xylane		<del></del>	4 <i></i>											
m and p xylene  Xylenes, total		-	ļ	<5.0 <5.0					T		<8.0 <5.0			

Compound	MW8-D 6/23/00	MFW8-D 9/22/00	MW8-D 11/30/00	MW9-D 7/26/99	MW9-D 9/9/99	MW9-D 11/1/99	MW9-D 3/2/00	MW9-D 6/23/00	MW9-D 9/29/00	MW9-D 11/30/00	MW1-WT 7/26/99	MW1-WT 9/9/99	11/1/99	MW1-WT 3/2/00
Acetone				<\$0	<50	<50					<50		<del></del> _	~~~~
Acrolein				<50	<50	<50	1		<del></del> -		<50			
Acrylonitrile	<u> </u>			<50	<50	<50			1		<50			<5
Benzene	<5.0	<5.0	<5.0	<5.0	<5.0		<5.0	<5.0	<5.0	<5.0	<5.0	<5.0	<5.0	<5.
Bromobenzene				<5.0	<5.0	<5.0					<5.0	<5.0	<5.0	<5.
Bromochloromethane		<u> </u>		<5.0	<5.0	<5.0					<5.0	<5.0	<5.0	45
Bromodichloromethane	ļ			<5.0	<5.0	<5.0					<5.0	<5.0	<5.0	45.
Bromoform		<u> </u>		<5.0	<50	<5.0					<5.0	<5.0	<5.0	<5.
Brome: - thane	<u> </u>			<10.0	<10.0	<10.0				<u> </u>	<10.0	<10.0	<10.0	<10.
n-butylbenzane				<10.0	<10.0						<10.0	<10.0	<10.0	<10.
sec-butylbenzene	<u> </u>			<10.0	<10.0	<10.0					<10.0	<10.0	<10.0	<10.
tert-butyženzene	<u> </u>			<10.0	<10.0	<10.0					<10.0	<10.0	<10.0	<10.
Carbon disulfide	<u> </u>			<10.0	<10.0	<10.0					<10.0	<10.0	<10.0	<10
Carbon totrachloride				<5.0	<5.0	<5.0					<5.0	<5.0	<5.0	<5
Ciriorobenzena				<5.0	<5.0	<5.0					<5.0	<5.0	<5.0	<5.
Chlorodibromomethane	L			<5.0	<5.0	<5.0					<5.0	<5.0	<5.0	<5.
Chioroethane				<10.0	<10.0	<10.0					<10.0	<10.0	<10.0	<10.
Chkroform		Ĭ _		<5.0	<5.0	<5.0					<5.0	<5.0	<5.0	<5.
Chloromethane				<10.0	<10.0	<10.0					<10.0	<10.0	<10.0	<10.
o-chiorotokiene				<5.0	<5.0	<5.0					<5.0	<5.0	<5.0	<b>&lt;</b> 5.
p-chlorotoluene	Ι			<5.0	<5.0	<5.0					<5.0	<5.0	<5.0	<5.
1,2-dibromo-3-chloropropane			L.							l				
1,2-dibromoethane										<del> </del>				<del></del>
Dibromomethane				<5.0	<5.0	<5.0				·	<5.0	<5.0	<5.0	<b>&lt;</b> 5.
1,2-dichlorobenzene	I			<5.0	<5.0	<5.0					<5.0	<5.0	<5.0	45.0
1,3-dichlorobenzene		I	·	<5.0	<5.0	<5.0				<b></b>	<5.0	<5.0	<5.0 <5.0	<5.1
p-dichlorobenzene				<5.0	<5.0	<5.0					<5.0 <5.0	<5.0 <5.0	<5.0 <5.0	<5.0 <5.0
Dichlorofluoromethane	Ī —	T				<del>-</del>				<del></del>	~3.0	13.0	<b>*3.0</b>	< 5.0
Dichlorodifluoromethane				<10.0	<10.0	<10.0					<10.0	<10.0		-40
1,2-dichloroethane		i		<5.0	<5.0	₹5.0					<5.0	<5.0	<10.0	<10.0
1,1-dichloroethane	<5.0	<5.0	<5.0	<5.0	<5.0	<5.0	<5.0	<5.0	<5.0	<5.0	<5.0		<5.0	<5.0
1,1-dichloroethane	<5.0	<5.0	<5.0	<5.0	<5.0	<5.0	<5.0	<5.0	<5.0	<5.0	<5.0	<5.0 <5.0	<5.0	<5.
cis-1,2-dichloroethene				<5.0	<5.0	<5.0			13.0	13.0	<5.0 <5.0	<5.0	<5.0	<5.0
rans-1,2-dichloroethene	1			<5.0	<5.0	<5.0					<5.0	<5.0	<5.0	<5.0
rans-1,4-dichloro-2-butone	T			<10.0	<10.0	<10.0					<10.0	<10.0	<5.0	< 5.0
di-Isopropylether	.)			1,0.0		1,0,0					< 10.0	<10.0	<10.0	<10,0
1,2-dichloropropane				<5.0	<5.0	₹5.0					<5.0	<5.0	45.0	
1,3-dichloropropane				<5.0	<5.0	<5.0					<5.0	<5.0	<5.0	<5.0
2,2-dichloropropane				<5.0	<5.0	<5.0					<5.0 <5.0	<5.0	<5.0	<5.0
1,1-dichloropropene				<5.0	<5.0	<5.0					<5.0	<b>45.0</b>	<5.0	<5.0
1,3-dichkropropene				<10.0	<10.0	<10.0					<10.0	<10.0	<5.0	<5.0
cis-1,3-dichloropropene				<5.0	<5.0	<5.0					<5.0	<5.0	<10.0	<10.0
rans-1,3-dichloropropene				<5.0	<5.0	<5.0					<5.0 <5.0	<5.0	<5.0 <5.0	<5.0 <5.0
trytenzene				<5.0	<5.0	<5.0				<del></del>	<5.0	<5.0	<5.0	
leptane												×3.0	· · · · · · · · · · · · · · · · · · ·	<5.0
2-hexanone				<5.0	<5.0	<5.0					<5.0	<5.0	<5.0	
iexachlorobutadiene				<10.0	<10.0	<10.0					<10.0	<10.0	<10.0	<5.0 <10.0
odomethane				<10.0	<10.0	<10.0					<10.0	<10.0	<10.0	<10.0 <10.0
sopropylbenzene				<10.0	<10.0	<10.0			<u> </u>		<10.0	<10.0	<10.0	
-isopropytoluene				<10.0	<10.0	<10.0					<10.0	<10.0	<10.0	<10.0
Aethylene chlorida				<5.0	<5.0	<5.0					<5.0	<5.0	<5.0	<6.0
Aethyl ethyl kelone				<10.0	<10.0	<10.0					<10.0	<10.0	<10.0	<10.0
ATBE											-10.0	*10.0		110.0
Asthyl isobutyl ketone		T												
-methyl-2-pentanone				<5.0	<5.0	<5.0					<5.0	<5.0	<5,0	25.0
etrahydrofuran	LI			<5.0	<5.0	<5.0					<5.0	<5.0	<5.0	<5.0
laphthalene														<b>-3.0</b>
propyBenzene				<10.0	<10.0	<10.0					<10.0	<10.0	<10.0	<10.0
tyrene			]	< 5.0	<5.0	<5.0					<5.0	<5.0	<5.0	<5.0
.1,1,2-letrachloroethane	T	7		<5.0	<5.0	<5.0					<50	<5.0	<5.0	<5.0
,1,2,2-tetrachloroethane				<5.0	<5.0	<5.0					<5.0	₹5.0	<5.0	<5.0
etrachicroethune				<5.0	<5.0	<5.0					<5.0 <5.0	<5.0	<5.0	<5.0 <5.0
oluane	<5.0	<5.0	<5.0	<5.0	<5.0	<5.0	<5.0	<5.0	<50	<5.0	<b>45.0</b>	<5.0	<5.0 <5.0	
,2,3-trichlorobenzene				<5.0	<5.0	<5.0		<del></del>			<5.0	<5.0 <5.0	<5.0 <5.0	<5.0
,2,4-trichlorobenzene				<5.0	<5.0	<5.0			<del>}</del>		<5.0	<5.0	<5.0	<5.0
,1,1-trichloroethane	<5.0	<5.0	<5.0	₹5.0	<50	<5.0	<5.0	<5.0	₹5.0	<5.0	<5.0	<5.0 <5.0		<5.0
1,2-trichloroethane				<5.0	<5.0	<5.0					<5.0	<5.0	<5.0	<5.0
richloroethene				<5.0	<5.0	<5.0					<5.0 <5.0	<5.0 <5.0	<5.0	<5.0
richloroflucromethane				<10.0	<10.0	<10.0					<10.0	<10.0	<5.0	<5.0
2,3-trichloropropane				<50	<5.0	<5.6			<del></del>		<5.0	<5.0	<10.0	<10.0
2,4-trimethylbenzone	1			<10.0	<10.0	<10.0					<10.0	<10.0	<5.0	<5.0
3,5-trimethylbenzene				<10.0	<10.0	<10.0					<10.0		<10.0	<10.0
nyl acelate	<del></del>			<10.0	<10.0	<10.0						<10.0	<10.0	<10.0
hyl acetate	}			- 10.0	-10.0]	- 10.0					<10.0	<10.0	<10.0	<10.0
nyl chloride	<del></del>			<10.0	<10.0	<10.0								
chloroethy Mnyl ether		+			-10.0	- 10.0		<b></b>		,	<10.0	<10.0	<10.0	<10.0
Tryl methacrylate							∤				l		I	
xylone		<del></del>	<del></del>	<10.0	<10.0	<10.0							I	
and p xylene			<u>-</u> -	<5.0							<10.0	<10.0	<10.0	<10.0
					<5.0	<5.0					<5.0	<5.0	< 5.0	<6.0
ienes, total				<5.0	<5.0	<5.0	- 1	- 1	1	· T	<5.0	<5.0	<5.0	<6.0

Maren	Compound	MW1-WT 6/23/00	MW1-WT 9/29/00	MW1-WT 11/30/60	MW2-WT 7/26/99	MWZ-WT \$/\$/89	MW2-WT	MW2-WT 3/2/00	MW2-WT 6/23/00	HW2-WT 9/29/00	MW2-WT	MW3-WT 7/26/99	MW3-WT
Marsell	Aceione		372000	1 (/30/00				3/2/00	023/00	3/23/00	11/30/00		<50
Security   Color   C								<del></del>					<50
Secondaries	<del></del>				\$	<5)	<50					<50	450
Second		<5.0	<5.0	<5.0				<5.0	<5.0	<5.0	<5.0		<5.0
Secretarian													<5.0
Secretary								<u> </u>		ļ	<u> </u>		<5.0
100   100	<del></del>			L				ļ					<5.0
Selegiphenses										ļ	ļ		<5.0
100   100								<b> </b>	<b></b>	<b></b>	ļ		
100   100								ļ	<del> </del>		<del> </del>		
Carbon devables								<del></del>		<del> </del>	<del></del>		
Carbon tendoride								<del></del>	<b></b>	}	<del></del>		<10.0
\$40   \$40										<del> </del>	<del> </del>		<5.0
Charles	Chiorobenzene												<5.0
Charbons	Chlorodibromomethane				<5.0					<b></b>			<5.0
Observations	Chiorce ihane				<10.0	<10.0	<10.0					<10.2	<10.0
	Chloroform				<5.0	<5.0	<5.0					<5.0	<5.0
Colorando   Colo					<10.0	<10.0	<10.0					<10.0	<10.0
13-80-many-hare				<u> </u>									<5.0
		<b> </b>			<5.0	<5.0	<5.0					<5.0	<5.0
September   Sept						<b></b>			<b></b>	ļ	ļ		
13-deltrotecures		<b> </b>		<b> </b>		<b></b>	<u></u>	<b> </b>	<b>!</b>	<b></b>	<b></b>	<u> </u>	ــــــــــــــــــــــــــــــــــــــ
13-del-informational   4-50   4-50   5-50   4-50				ļi				<b></b>	<del></del>	ļ	ļ		<5.0
Section   Sect		<del></del>						<del> </del>	<del> </del>	<del> </del>	<del> </del>		<5.0
Debreimbursomburs		<del> </del>		<del> </del>				<del> </del>	<del> </del>	<del> </del>	<del> </del>		<5.0 <5.0
100   100				<del> </del>	7.0			<del> </del>	<del>                                     </del>		<del> </del>	₹3.0	
12-doll-brownem	· · · · · · · · · · · · · · · · · · ·			<b> </b>	<10.0	<10.0	<10.0		<del></del>		<del> </del>	€10 Å	<10.0
11-6de/browhene	1,2-dichlorcethane									<del> </del>	<del> </del>		<5.0
11-dehitopropere	1,1-dichloroethane	<5.0	<5.0	<5.0	<5.0	<5.0	<5.0	<5.0	<5.0	<5.0	<5.0		45.
Team   1,50 district property   1,50 distric	1,1-dichloroethene	<5.0	<5.0	<5.0	<5.0	<5.0	<5.0	<5.0	<5.0	<5.0	<5.0		7.90
Table   Adultive Delatine					<5.0	<5.0	<5.0					<5.0	<5.0
dispropry/defer												<5.0	<5.0
12-delitopropopee					<10.0	<10.0	<10.0					<10.0	<10.0
13.49 https://proposes								ļ	ļ				
22-dichtoproposes											<b> </b> -		<50
1.1.dishbropropense									<u> </u>	ļ			45.0
1.3-dechtoproposes				<u> </u>				ļ	<del></del>		ļ		<5.0
dels   3.4-delshoropropens								<del></del>		<del> </del>	<del></del>		
								<del> </del>	<del></del>		<del> </del>		<5.0
Ellybename										<del> </del>			<5.0
Heptene	Ethylbenzene				<5.0	<5.0					<u> </u>		<5.0
Managhtopicalative	Heptane												<del>-</del> -
Decomphisms	·				<5.0							<5.0	<5.0
IsopropyBenzane												<10.0	<10.0
P-1000000000000000000000000000000000000													<10.0
Methylethore										<u> </u>	ļ		<10.0
Methyl ethyl ethons										ļ <u> </u>	<u> </u>		<10.0
MBE										<del> </del>			<5.0
Methyl lockuryl ketone					*10.0	*10.0	¥10.0		<del> </del>	<del></del> -	<del> </del>	<10.0	<10.0
4-methyl-2-pentanone								<del></del>	<del>                                     </del>	<del> </del>	<del> </del>	<del></del>	<del> </del> -
Test alty-forus					<5.0	<5.0	<5.0	<del></del>	<del></del>	<del></del>	†	<50	<5.0
Naphthales	Tetrahyúrofuran								L				<5.0
Styrene	Naphthalene												
1,1,1,2-lev/achloroethane	n-propythenzene			ļ								<10.0	<10.0
1,1,2 2-tar schloroethane													<5.0
Tetrachloroethene									<b> </b>				<5.0
Tobset											ļ		<5.0
1,2,3-tichlorobenzene	<del></del>												<5.0
1.2.4 trichlorobenzene		<b>*5.U</b>	*3.0	<5.0				<5.0	<5.0	<5.0	<5.0		<5.0
1,1,1-trichtoroethane								<del></del>	<del></del>	<del></del>	<del> </del>		<5.0
1,1,2-tichloroethane	1,1,1-trichloroethane	<5.0	<5.0	<5.0				<4.0	ck n	28 n	£60		<5.0
Trichloroethene         <50	1,1,2-trichloroethane				•				<u></u>		<del></del>		<5.0
Trichkroftvorome*tane	Trichloroethene								l		<del>                                     </del>		₹5.0
1,2,3 trinshpropropers	Trichlorofluoromethane										<del>                                     </del>		<10.0
1.2.4 trimetry/benzerie	1,2,3-trichtoropropene	I											<5.0
1,3,5 trimetry benzene	1,2,4-trimethy/benzerie					~					L		<10.0
Vinyl acetale         < 10.0	1,3,5-trimethylbenzene												<10.0
Etyl acetals	Vinyl acetata	]	[		<10.0	<10.0	<10.0						<10.0
Critical triple	Ethyl acelate				]								
C-chicarolityNdry  e89er	Virryl chloride				<10.0	<10.0	<10.0					<10.0	<10.0
- xylane	2-chioroethyMnyl ether												
m and p sylene <50 <50 <50 <50 <50 <50 <50 <50 <50 <50	Ethyl methacrylate												
	o-xylene												<10.0
	m and p xylene Xylenes, lotal								ļ	ļ	ļ		

									·			
Compound	MW3-WT	MW3-WT	MV/3-WT	MW3-WT	MW3-WT	MW4-WT	MW4-WT	MW4-WT	MW4-WT	MW4-WT	MW4-WT	MMEMI
	11/1/99	3/2/09	6/23/00	9/29/00	11/30/00	7/26/99	9/9/99	11/1/99	3/2/00	6/23/00	9/29/00	11/30/00
Acetons	<50 <50					<50	<50	<50				
Acrolein Acrylonitrile	450				·	<50 <50	<50 <50	<50 <50				
Benzene	<5.0	<5.0	<5,0	<5.0	<5.0	<5.0	45.0	<5.0	<5.0	<5.0	<5.0	<5.0
Bromobenzane	<5.0					<5.0	<5.0	<5.0				
Bromochloromethane	<5.0					<5.0	<5.0	<5.0				
Bromodichloromethane	<5.0					<5.0	<5.0	<5.0				
Bromoform	<5.0					<5.0	<5.0	<5.0				
Bromomethane	<10.0					<10.0	<10.0	<10.0				
n-buty/benzene sec-buty/benzene	<10.0 <10.0					<10.0 <10.0	<10.0 <10.0	<10.0 <10.0				
iert-butybanzene	<10.0					<10.0	<10.0	<10.0				
Carbon disulfide	<10.0					<10.0	<10.0	<10.0				
Carbon tetrachloride	<5.0					<5.0	<5.0	<5.0				
Chlorobenzene	<5.0					₹ <b>3.</b> 0	<5.0	<5.0				
Chlorodibromomethane	<5.0					<5.0	<b>45.0</b>	<5.0				
Chiorcethane	<10.0		ļ			<10.0	<10.0	<10.0				
Chloroform	<5.0					6.7	<5,0	<5.0				
Chloromethane	<10.0					-10.0	<10.0	<10.0				<del></del>
p-chicrotoluene	<5.0 <5.0					<5.0 <5.0	<5.0 <5.0	<5.0 <5.0				
1,2-dibromo-3-chloropropane		<del> </del>					45.0	<b>10.0</b>				
1,2-dt.romoethane												
Dibromometriane	<5.0					₹5.0	<5.0	<5.0				
1,2-dichlorobenzene	<5.0					<5.0	<5.0	<5.0				
1,3-dichlorobenzene	<5.0					<5.0	<5.0	<5.0				
p-dichlorobenzene	<5.0					<5.0	<5.0	<5.0				
Dichloroft.crome*hane		<b> </b>										
Dichlorodifluoromethane 1,2-dichloroethane	<10.0 <5.0					<10.0	<10.0 <5.0	<10.0 <5.0				
1.1-dichloroethane	8.1	<5.0	<5.0		6.50	<5.0 <5.0	<5.0	<5.0	<5.0	<5.0	45.0	<5.0
1,1-dichloroethens	<5.0	<5.0	<5.0	<5.0	<5.0	<del> </del>	<b>&lt;5.0</b>	<5.0	<5.0	<b>45.0</b>	<b>45.0</b>	45.0
ds-1,2-dichloroethene	<5.0					<5.0	<5.0	<5.0				
trans-1,2-dichloroethene	<5.0					<5.0	<5.0	<5.0				
trans-1,4-dictioro-2-butene	<10.0					<10.0	<100	<10.0				
di-isopropylether												
1,2-dichloropropane	<5.0					<50	<5.0	<5.0				
1,3-dichioropropane	<5.0		ļ			<5.0	<5.0	<50				
2.2-dichloropropane	<5.0 <5.0		ļ			<5.0	<5.0	45.0				
1,1-dichloropropene	<10.0		<del></del>			<5.0 <10.0	<5.0 <10.0	<5.0 <10.0			<u> </u>	
cit-1,3-dichloropropene	<5.0					15.0	<5.0	<5.0			<del></del>	
trans-1,3-dichloropropene	<5.0		<del> </del>			₹50	<5.0	<5.0			·	
Eftylbenzene	<5.0	·				<5.0	<5.0	<5.0		· · · · ·		
Heptane												
2-haxanone	<5.0					<5.0	<5.0	<5.0				
Hexachlorobutadiene	<10.0		<b></b>	L		<10.0	<10.0	<10.0				
lodomethane	<10.0 <10.0					<10.0	<10.0 <10.0	<10.0				
laopropy/benzene p-isopropy/liotuene	<10.0		<del> </del>	<b></b>	<del></del>	<10.0 <10.0	<10.0	<10.0 <10.0				
Methylene chloride	<5.0		<del></del>			<5.0	<5.0	<5.0				
Methyl ethyl kelona	<10.0		<u> </u>	<b> </b>		<10.0	<10.0	<10.0				
MTBE			I									
Mathyi isobutyi kelone	ļ											
4-methyl-2-pentanone	<5.0		ļ			<5.0						
Tetratrydrofuran Naphthalene	<5.0	ļ	ļ	<b> </b> -		<5.0	<5.0	<5.0	ļ			<b></b>
n-propylbenzene	<10.0	<del></del>		<del> </del>		<10.0	<100	<10.0	<del> </del>			<del> </del>
Styrene	<5.0		<del> </del>		<del></del>	<5.0						
1,1,1,2-tatrachlorgethane	<5.0				····	<5.0					<b></b>	<del> </del>
1,1,2,2-letrarhibroethane	<50					<5.0					l	
Tetrachioroethene	<5.0					<5.0						
Toluene	<5.0		<5.0	<5.0	<5.0	<5.0	<5.0			<5.0	<5.0	<5.0
1,2,3-trichlorobenzene	<5.0		ļ	ļ		<5.0						
1,2,4-tichlorobenzene	<5.0 35.6					<5.0				ļ	ļ <u>.</u>	
1,1,1-\$ ichioroethane 1,1,2-trichioroethane	J5.0 <5.0	18	19.7	34.1	22.6	<5.0 <5.0				<5.0	<5.0	<5.0
Trichloroethene	450			<del> </del>		<5.0 <5.0					<del> </del>	
Trichlorofluoromethane	<10.0					<10.0	<10.0			<del></del>	<del> </del>	<del></del>
1,2,3-trichkropropane	<5.0		<del></del> -	······		<5.0	<5.0			<del> </del>	<del> </del> -	r
1,2,4-trimethylbenzene	<10.0					<10.0				<del></del>	<del> </del>	
1,3,5-trimethy/benzene	<b>410.0</b>	<u> </u>				<10.0					1	
Virryl acetato	<10.0					<10.0	<100	<10.0				
Ethyl acetate												
Vinyl chloride	<10.0					<10.0	<10.0	<10.0	L	<u> </u>		
2-chlorcethylvinyl ather	ļ											
Ethyl methacrylate	240.5							<b></b>	<b> </b>		ļ	
o-xylene m and p xylene	<10.0 <5.0					<10.0 <5.0		<5.0		<del> </del>	<b> </b>	ļl
m and p xysene Xylenes, fotal	<50					<5.0				ļ		
ngan wa, 'sser	130	ن ــــــــــــــــــــــــــــــــــــ	I			< 5.0	₹5.0	49.0	·	<u></u>	I	i

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Compound	IW-WI	MW9-WT	MW9-WT	MW9-WT	MW9-WT	1217.10	MW9-WT	LW-1	LW-1	LW-1	LW-1	LW-1	LW-1
Compound	7/26/99	9.9.99	11/1/99	3/2/00	6/23/00	M14/9-W/T 3/29/00	11/30/00	6/17/98	1/24/98	10/26/98	7/26/99	9/9/93	11/1/99
Acetone	<b>450</b>	<50	<50	3200	023.00	37500	1113400	<20	<20	<20	<50	₹50	<50
Acrolein	<50	<50	<b>450</b>		<u> </u>				<10	<50	<50	<50	<50
Azrylonitrile	<50	<50	<b>-</b> 50						<10	<10	<50	<50	<50
Benzene	<5.0	<5.0	<5.0	<5.0	<5.0	<5.0	<5.0	1.4	<0.5	<0.5	<5.0	<5.0	<5.0
Bromobenzene	<5.0	<5.0	<5.0					<1.0	<1.0	<1.0	<5.0	<5.0	<5.0
Bromochloromethane Bromodichloromethane	<5.0 <5.0	<5.0 <5.0	<5.0 <5.0					<1.0 <1.0	<1.0 <1.0	<1.0 <1.0	<5.0 <5.0	<5.0 <5.0	<5.0 <5.8
Bromoform	45.0	45.0	<5.0					<2.0	<20	<2.0	<5.0	<5.0	<5.0
Bromomethane	<10.0	<10.0	<10.0					c4.0	<4.0	4.0	<10.0	<10.0	<10.0
n-buty/benzene	<10.0	<10.0	<10.0					<1.0	<1.0	<1.0	<10.0	<10.0	<10.0
sec-buty/benzene	<10.0	<10.0	<10.0					<1.0	<1.0	<1.0	<10.0	<10.0	<10.0
tert-butytsenzene	<10.0	<10.0	<10.0					<1.0	<1.0	<1.0	<10.0	<10.0	<10.0
Carbon disulfide	<10.0	<10.0	<10.0						<1.0	<1.0	<100	<100	<10.0
Carbon tetrachloride	<5.0	<5.0	<5.0	ļ				<1.0	<1.0	<0.3	<5.0	<5.0	<5.0
Chlorobenzene Chlorodibromomethane	<5.0 <5.0	<5.0 <5.0	<5.0 <5.0	<del></del>				<1.0 <1.0	<1.0 <1.0	<1.0 <1.0	<5.0 <5.0	<5.0 <5.0	<5.0 <5.0
Chloroethane	<10.0	<10.0	<10.0	<b></b>	<u> </u>			<4.0	<4.0	4.0	<10.0	<10.0	<10.0
Chlorotom	45.0	<5.0	<5.0	<del> </del>				<1.0	<1.0	<1.0	<5.0	<b>45.0</b>	₹5.0
Chlororrathane	<10.0	<10.0	<10.0					<1.0	<1.0	<1.0	<10.0	<10.0	<10.0
a-chlorotoluene	<5.0	<5.0	<5.0		<del> </del>	<del></del>		<1.0	<1.0	<1.0	<5.0	<5.0	<5.0
p-chlorotokusne	<5.0	<5.0	<5.0					<1.0	<1.0	<1.0	<5.0	<50	<\$.0
1,2-dibromo-3-chloropropane								<10	<10	<1.0			
1,2-dibromoethane			ļ	ļ				<10	<10	<10		1	
Dibromomethane	<5.0	<5.0	<50	ļ	<b></b>			<1.0	<1.0	<1.0	<5.0	<5.0	<5.3
1,2-dichlorobenzene 1,3-dichlorobenzene	<5.0 <5.0	<5.0 <5.0	<5.0 <5.0	<del> </del>	<b></b>			<1.0	<1.0	<1.0	<5.0	<5.0 <5.0	<5.0 <5.0
p-dichlorobenzene	<5.0	<5.0 <5.0	<5.0 <5.0	<del> </del>	<b></b>	<b></b>		<1.0 <1.0	<1.0 <1.0	<1.0	<5.0 <5.0	<u>4.0</u>	<5.0 <5.0
Dichlorofluoromethane		<del></del>	1 3.0	<del> </del>		<del></del>	<del></del>	1.0	<10	43.0	1	90	19.0
Dichlorodifluoromathane	<10.0	<10.0	<100	l	<del> </del>			<3.0	<3.0	43.0	<10.0	<10.0	<10.0
1,2-dichloroe/hane	<5.0	<5.0	<5.0					<1.0	<1.0	<1.0	<5.0	<5.0	<5.6
1,1-dichloroethane	<50	8.30	5.3	13.2	8.5	10.4	6.3	1.3	<1.0	1.3	<5.0	<b>&lt;</b> 5.0	4.
1,1-dichloruethene	<50	<5.0	<50	<5.0	<5.0	<50	<5,d	42.0	<20	42.3	<5.0	<5.0	<5.0
cis-1,2-dichloroethene	<5.0	<5.0	<5.0	<b></b>	ļ			<1.0	<1.0	<1.0	<5.0	<5.0	<5.0
trans-1,2-dichloroethene	45.0	<5.0	<5.0	<b></b>	ļ	ļ		<1.0	<1.0	<1.0	<5.0	₹5.0	<5.0
trans-1,4-dichloro-2-butene di-isopropylether	<10.0	<10.0	<100	<u> </u>	ļ			<1.0	<10 <1.0	<10	<10.0	<10.0	<10.0
1,2-dichloropropane	40	<5.0	<50	<del> </del>	<del> </del>			<1.0	<1.0	<1.0 <1.0	<5.0	<50	<5.0
1,3-dichloropropane	45.0	<5.0	<5.0		<del></del>			<1.0	<1.0	<1.0	<5.0	<b>45.0</b>	<5.0
2,2-dichloropropane	<5.0		<5.0	<del></del>				<1.0	<1.0	<1.0	<5.0	<5.0	<5.0
1,1-dichloropropene	<5.0	<5.0	<5.0					<1.0	<1.0	<1.0	<50	<5.0	<5.0
1,3-dichloropropene	<10.0		<10.0								<10.0	<10.0	<10.0
cis-1,3-dichloropropene	<5.0	<5.0	<50					<1.0	<1.0	<1.0	<5.0	<5.0	<5.0
trans-1,3-dichloropropene Ethylbenzene	45.0 45.0	<5.0 <5.0	<5.0 <5.0	ļ	<u> </u>			<1.0	<1.0	<1.0	<5.0	<5.0	<5.0
Heptane	43.0		*3.0	<del> </del>	<del></del>			<1.0	<1.0 <1.0	<1.0	<50	<5.0	<5.0
2-hexanone	40	<5.0	<5.0	<del> </del>	<del> </del>				<10	<10	<5.0	<5.0	<5.0
Horachiorobutadiene	<100	<10.0	<10.0		<del> </del>			<5.0	₹5.0	<5.0	<10.0	<10.0	<10.0
lodomethane	<10.0	<10.0	<10.0						<10	<10	<10.0	<10.0	<10.0
isopropy/benzene	<10.0	<10.0	<100					<1.0	<1.0	<1.0	<10.0	<10.0	<10.0
p isopropylloluene	<10.0	<10.0	<10.0		ļ			<1.0	<1.0	<1.0	<10.0	<10.0	<10.0
Methylene chloride Methyl ethyl ketone	<5.0 <10.0	<50 <100	<5.↑ <10 U		ļ			<10	<10	<5.0	<5.0	<5.0	<5.0
MTBE	1100		·····	<del> </del>	<del> </del>	<del> </del>		<10 <1.0	<10 <1.6	<10 <1.0	<10.0	<10.0	<10.0
Methyl isobutyl ketone	l			<del>                                     </del>	<del> </del>		<u> </u>	<u></u>	<10		<u> </u>		
4 methyl-2-pentanone	<50	<50	<5.0			•				<u></u>	<5.0	<5.0	<5.0
Tetrahydrofuran	<50	<50	<50						<50			<50	<5.0
Maphithalene				ļ	ļ			<5.0					
n-propybenzere	<10 0 <5.0	<10.0 <5.0	<10.0 <6.0		<b> </b>		<b></b> _	<1.0	<1.0			<10.0	<10.0
Styrene 1,1,1,2 tetrachloroethane	<5.0 <5.0		<5.0		<del></del>			<1.0 <1.0	<1.0 <1.0			<5.0	<5.0
1,1,2,2-tetrachioroethane	<5.0	<5.0	₹5.0		·			<1.0	<1.0			<5.0	<5.0
Tetrachkmetiene	<5.0	15.0	45.0					<1.0	<10			<5.0	45.0
Totuena	<50	<50	<5.0	<5.0	<5.0	<50	<50	1.2	<1.0			<5.0	<5.D
1,2,3 trichlorobenzene	<b>&lt;</b> 50		<5.0	\$ <del></del>		L		<5.0	<5.0	<5.0		<5.0	<5.0
1,2,4 Inchlorobenzene	<50		<5.0	•				< 5.0	₹60		<5.0	₹5.0	<5.0
1,1,1-tichlorcethane		<50	9.5	41.3	10.5	11.3	15.3	<1.0	<1.0			<5.0	45.0
1,1,2 trichloroetrane	<5.0	<50	<5.0	·	ļ			<1.0	<1.0			<6.0	<5.0
Trichlorcethene Trichlorofluoromethane	<5 C	<50 <100	<5.0 <10.0	<del> </del>	·			<1.0 <40	<1.0		<del></del>	<5.0	<50
1.2,3 trichicropropane	<5.0		<5.0	<del> </del>				<1.0	<1.0	<4.0 <1.0		<10.0 <5.0	<10.0
1,2,4 timetrytenzene	<10.0	<10.0	<10.0	<b> </b>				<1.0	<1.0			<10.0	<5.0 <10.0
1,3,5 timetry/benzene	<10.0	<10.0	<10.0	<del> </del>				<1.0	<del>&lt;1.0</del>			<10.0	<10.0
Vinyl anetate	<10.0	<10.0	<10.0	<del> </del>				<del></del>	<b>410</b>			₹10.0	
Ethyl acetate									l	l			
Vinyl chloride	<10.0	<10.0	<10.0					<1.0	<1.0	<1.0	<10.0	<10.0	<10.0
2-chlorcethylvinyl ether	•			ļ									
Ethyl mathacrylsia				ļ							ļ		
o-sylene m end a rolena	<10.0 <5.0	<10.0 <5.0	<10.0 <5.0		ļ		ļ		ļ	ļ	<10.0	<10.0	<10.0
m and p xylone Xylenes, lotal	<50 <50		<50		·			<30	<30	43.0	<50	<6.0	<5.0
		الارد-		L	ز ــــــــــــــــــــــــــــــــــــ	نسيرسيب		*30	·	1 430	<5.0	<\$.0	<5.0

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Compound	IW9-WT	MWS-WT	MW9-WT	MWS-WT	MW9-WT	4040 W/T	MW9-WT	LW-1	LW-1	LW-1	LW-1	LW-1	LW-1
Compound	1/26/99	9/9/99	11/1/99	3/2/00	6/23/06	9/29/00	11/30/00	6/17/98	1/24/98	10/26/98	1/26/99	9/9/99	11/1/99
Acelone	<50	<50	<50					<20	<20	<20	<50	< <b>50</b>	<50
Acrolein	<\$0	<50	<b>450</b>						<10	<b>&lt;</b> 50	<50	<50	<50
Acrylonárile	<50	<50	<50						<10	<10	₹50	<50	<50
Senzene	<5.0	<5.0	<5.0	<5.0	<5.0	<5.0	<\$.0	1.4	<0.5	<0.5	<5.0	<5.0 <5.0	<5.0
Bromobenzene Bromochloromethane	<5.0 <5.0	<5.0 <5.0	<5.0 <5.0					<1.0 <1.0	<1.0 <1.0	<1.0 <1.0	<5.0 <5.0	<5.0	<5.0
Bromodichlorome@nane	450	<b>45.0</b>	<5.0					<1.0	<1.0	<1.0	<5.0	<5.0	<5.0
Promotorm	<5.0	<5.0	<5.0					<2.0	<2.0	<2.0	<5.0	<5.0	<5.0
Bromomethane	<10.0	<10.0	<10.0					<4.0	44.0	<4.0	<10.0	<10.0	<10.0
n-buly/benzene	<10.0	<10.0	<10.0					<1.0	<1.0	<1.0	<10.0	<10.0	<10.0
sec-butythenzene	<10.0		<10.0					<1.0	<1.0	<1.0	<10.0	<10.0	<10.0
tert-butythenzene	<10.0		<10.0	<b></b>				41.0	<1.0	<1.0	<10.0	<10.0	<10.0
Carbon disulfide Carbon tetrachloride	<10.0 <5.0	<10.0 <5.0	<10.0 <5.0		ļ				<1.0 <1.0	<1.0	<10.0 <5.0	<10 0 <5.0	<10.0 <5.0
Chlorobenzene	<5.0	<5.0	< <u>&lt;5.0</u>		<del> </del>			<1.0 <1.0	<1.0	<1.0	<5.0 <5.0	45.0	<5.0
Chlorodibromomethane	40		<5.0		<del> </del>			<1.0	<1.0	<1.0	<5.0	<50	<5.0
Chloroethane	<10.0		<10.0		·			41.0	44.0	<4.0	<10.0	<10.0	<10.0
Chloroform	<5.0	<5.0	<5.0					<1.0	<1.0	<1.0	<5.0	<5.0	<5.0
Chlororrathane	<10.0	<10.0	<10.0					<1.0	<1.0	<1.0	<10.0	<10.0	<10.0
o-chlorotoluene	<5.0		<50					<1.0	<1.0	<1.0	<5.0	<5.0	< 5.0
p-chlorotolusne	<5.0	<5.0	<5.0		<u> </u>			<1.0	<1.0	<1.0	<5.0	<5.0	<5.0
1,2-dibromo-3-chioroprogane	ļ	<b></b>	L	ļ				<10	<10	<1.0	ļ	<u>                                     </u>	
1,2-dibromoethane Dibromomethane	<5.0	<5.0	<5.0	<del> </del>	<del> </del>			<10	<1.0	<10	<b></b>	<5.0	<5.3
1,2 dichlorobenzerie	<5.0		<5.0 <5.0	<del> </del>	<del> </del>			<1.0 <1.0	<1.0 <1.0	<1.0 <1.0	<5.0 <5.0	<5.0 <5.0	<5.0 <5.0
1,3-dichlorobenzene	40		<5.0	t	<del> </del>			<1.0	<1.0	<1.0	₹5.0	<5.0	<5.0
p-dichlorobenzene	45.0		<5.0	<del></del>	<b></b>			<1.0	<1.0	<1.0	<5.0	30	45.0
Dichlorofluoromethane	1	<del></del>		<b> </b>	<b>†</b>				<10	<3.0	<u>                                     </u>	<del>-</del>	
Dichlorodifluoromathane	<10.0	<10.0	<10.0					<3.0	<3.0	<3.0	<10.0	<10.0	<10.0
1,2-dichloroe/hane	<5.0		<5.0					<1.0	<1.0	<1.0	<5.0	<5.0	< 5.0
1,1-dichloroethane	450	8.30	5.3	13.2	8.5	10.4	6.3	1.3	<1.0	1.3	<5.0	<5.0	45.
1,1-dichloruethene	<50		<50	45.0	<5.0	<50	<5.J	<2.0	<2.0	<2.0	<5.0	<5.0	<5.0
cis-1,2-dichloroethene trans-1,2-dichloroethene	<5.0 <5.0		<5.0 <5.0	<del> </del>	<del> </del>			<1.0 <1.0	<1.0 <1.0	<1.0 <1.0	<5.0	<5.0 <5.0	4.0 45.0
tans-1,4-dichloro-2-butene	<10.0		<100	<del> </del>		<del></del>		<1.0	<10	<10	<5.0 <10.0	<10.0	<10.0
di isopropylether		1	[—— <del></del>	<del> </del>			~	*****	<1.0	<1.0	×10.0		- 100
1,2-dichloropropane	40	<5.0	450	<b></b>	<del> </del>			<1.0	<1.0	<1.0	<5.0	<5.0	<5.0
1,3-dichloropropane	<5.0	<50	<5.0					<1.0	<1.0	<1.0	<5.0	<5.0	<5.0
2,2-6chloroprosane	<5.0		<5.0					<1.0	<1.0	<1.0	<5.0	<5.0	<5.0
1,1-dichloropropene	<5.0		₹5.0		ļ			<1.0	<1.0	<1.0	<50	<5.0	<5.0
1,3-dichkropropena	<10.0 <5.0		<10.0 <5.0		ļ						<10.0	<100	<10.0
cis-1,3-dichloropropene trans-1,3-dichloropropene	450		<50 <50		<del> </del>			<1.0 <1.0	<1.0 <1.0	<1.0 <1.0	<5.0 <5.0	<5.0 <5.0	<5.0
Ethylbenzene	40		<5.0	<del> </del>	<del> </del>			11.0	<1.0	<1.0	<50	<5.0	<5.0 <5.0
Heptane		<u> </u>							<1.0		1		
2-hexanone	<50	<50	<5.0						<10	<10	<5.0	<5.0	<5.0
Herachiorobutadiene	<100		<10.0	<del></del>				<5.0	<5.0	<5.0	<10.0	<10.0	<10.0
lodomethane	<10.0		<10.0	<u> </u>					<10	<10	<10.0	<10.0	<10.0
isopropytherzene	<10.0 <10.0	<del></del>	<10.0 <10.0		<del> </del>			<1.0	<1.0	<1.0	×10.0	<10.0	<10.0
p-isopropyllokuene Multhylene chloride	<50		€5,7		<del> </del>			<1.0 <10	<1.0 <10	<1.0 <5.0	<10.0 <5.0	<10.0 <5.0	<10.0 <5.0
Methyl ethyl ketone	<10.0		<10 U		<del> </del>			<10	<10			<10.0	<10.0
NTBE		I		<u> </u>	1		<del></del>	<1.0	₹1.6	<1.0	1		
Methyl isobutyl ketone		<u> </u>							<10	<b>≤10</b>			
4 methyl-2-pentanone	<50	<50	<5.0	<u> </u>	ļ						<5.0	<5.0	<5.0
Tetrahydrofuran	<50	<50	<50	<b> </b>	<del> </del> -	<u> </u>			<50	<50	<50	<5.0	<5.0
Naphitralene n-propythenzene	<10.0	<100	<10.0	<del> </del> -	<del> </del> -	ļ		<5.0 <1.0	<5.0 <1.0			<u></u>	
Styrene	<5.0		₹5.0		<del> </del>	<del></del>	<del></del>	<1.0	<1.0			<10.0 <5.0	<10.0 <5.0
1,1,1,2 tetrachioroethane	<50		<50		<u> </u>			<1.0	<1.0			<5.0	<5.0
1,1,2,2-tet achioroethane	<5.0	<50	<5.0					<1.0	<1.0			₹5.0	<5.0
Tetrachlomethene	<50		<5.0					<1.0	<10	<1.0	<5.0	<5.0	<60
Toluena	<50		<5.0		<50	<50	<50	1.2	<1.0				
1,2,3 trichlorobenzene	<b>450</b>		<5.0	<del></del>	<u> </u>	ļ		<5.0	<5.0	< 5.0			<5.0
1,2,4-trichlorobenzene	<50	<5.0 <5.0	<5.0 9.5	41.3	15.6		15.3	<5.0 <1.0	<50			<5.0	<5.0
1,1,2 tichloroetiane	<5.0		<5.0		11.5	11.3	10.3	₹1.0 ₹1.0	<1.0 <1.0			<5.0 <6.0	<5.0 <5.0
Trichlorcethene	√5 C		<5.0		<del>                                     </del>			<1.0	<1.0				<5.0 <5.0
Trichlorofluoromethane	<10.0		<10.0					<40	<4.0				<10.0
1,2,3 trichloropropane	<5.0		< 5.0					<1.0	<1.0				45.0
1,2,4 trimethy/tenzene	<100		<10.0					<1.0	<1.0	<del></del>	<100		<10.0
1,3,5 trimetrytenzene	<100		<10.0					<1.0	<1.0			<10.0	<10.0
Vinyl acetate	<10.0	<10.0	<10.0	ļ	ļ				<10	<10	<10,0	<10.0	<10.0
Ethyl acotate				ļ	ļ <u>.</u>	<del></del>		ļ	ļ	ļ			L
Vinyl chloride	<10.0	<10.0	<10.C	<del> </del>	ļ	<del> </del>		<1.0	<1.0	<1.0	<10.0	<10.0	<10.0
								i	1	ı	1		ı
2-chlorcethylvinyl ether					<del> </del>				<del></del>	<del> </del>	·	<del></del>	
2-chlorcethylvinyl ether Ethyl methacrylsia		<10.0	<10.0								¢10.0	z10.0	400
2-chilorcethylvinyl ether	<10.0 <5.0		<10.0 <5.0								<10.0 <5.0	<10.0 <5.0	

Compound	LW-1	LW-1	LW-1	LW-1	LW-3	LW3	LW-3	LW-3	LW-3	LW-3	SPRING	SPRING	SP70NG	SPRING
·	3/2/00	67350	9/29/00	11/30/09	7/26/99	9.9.99	11/1/99	6/23/00	9/79/00	11/30/00	7/26/99	9/9/99	1171/99	3/2/00
Acetorie					<50	<50	<50				<50	<50	<50	
Acrolein						<50	<50				<50	88	<50	
Acrylonitrie Benzene	<5.0	<5.0	<5.0	<5.0	<50 <5.0	<50 <5.0	<50 <5.0	<50	<€ 3	<5.0	<50	<50 <5.0	<50 <5.0	<5.0
Bromobenzene				9.0	45.0	45.0	<5.0	(3.0	6.3	9.0	<5.0	<5.0	45.0	- 45.0
Bromochkromethane		<b></b>	<u> </u>		€0	<5.0	<50				45.0	<5.0	₹5.0	
Bromodohkromethare					0.0	<5.0	<5.0				<5.0	<5.0	<5.0	
Bromolorm					<b>450</b>	<5.0	<5.0				<b>&lt;</b> 5.0	<5.0	<5.0	
Bromomethane					<10.0	<10.0	<10.0				<10.0	<10.0	<10.0	
n-butybenzere					<10.0	<10.0	4177				<10.0	<10.0	<10.0	
sec-butytoenzene tert-butytoenzene	<b></b>	<b> </b>			<10.0 <10.0	<10.0 <10.0	<10.0 <10.0				<10.0	<10.0 <10.0	<10.0 <10.0	
Carbon disulfide	-				<10.0	¢10.0	<10.0				<10.0 <10.0	<19.0	<10.0	
Carbon tetrachioride	<del></del>				<50	<5.0	<5.0				<5.0	<5.0	45.0	
Chloroberzene					4.0	<5.0	<5.0				<5.0	<5.0	<5.0	
Chlorod ,romamethane					<5.0	<5.0	<5.0				<2.0	<5.0	<5.0	
Chiorathica	ļ	<u> </u>	ļ		<10.0	<10.0	<10.0				<10.0	<10.0	<10.0	
Chloroform Chlorometiane	ļ	<del> </del>			<50	<5.0	<5.0				<5.0	<5.0		
o-chiorotoluene	<del> </del>				< 10.0	<10.0 <5.0	< 10.0 < 5.0				<10.0 <5.0	<10.0 <5.0	<10.0 <5.0	
p-chiorotoluene		<b></b>	<del>                                     </del>		4.0	<5.0	45.0				<5.0	<b>45.0</b>	<50	
1.2-dikromo-3-chicropropane		ì	1		7:3		—— <u>"</u>							
1,2-ditromoethure														
Obromomethane		ļ			<5.0	<5.0	<5.0				<5.0	<5.0	<5.0	
1,2-dichlorobenzene	ļ	<b> </b>	ļ		<b>4.0</b>	<50	<5.0				<5.0	< 5.0	<5.0	
1,3-dichlorobenzene	<del> </del>	<del> </del>			<5.0 <5.0	<5.0	<5.0 <5.0		<b></b>		<5.0	<5.0 <5.0	<5.0 <5.0	
Dichlarofluoromethane	<del>                                     </del>	<del> </del>	<del> </del>	<del></del> -	9.0	49.0			<del></del>	ļ	<5.0	<5.0	9.0	
Dichlorodifucromethane	<b></b>	<b> </b>	<b> </b>		<100	<10.0	<10.0				<10.0	<10.0	<10.0	
1.2-dichlorosthane					<50	<5.0	<5.0				<5.0	<5.0	45.0	
1,1-dichloroethane	<5.0	<del></del>		<5.0	<5.0	<5.0	<5.0	<5.0	<5.0	45.0	<5.0	<5.0	<5.0	<\$.
1,1-dichloroethens	45.0	<5.0	<5.0	<5.0	<50	45.0	<50		<5.0	<50	45.0	<5.0	45.0	<b>45</b> .
cis-1,2-dichicrosthene trans-1,2-dichicrosthene	<del> </del>	<del> </del>	<b> </b>		<50	<5.0 <5.0	<50				<5.0	<50	<5.0	
trans-1,4-dichloro-2-bulane	<del> </del>		<del> </del>		<5.0 <10.0	<10.0	<5.0 <10.0	ļ			<50 <100	<5.0 <10.0	<5.0 -10.0	
di-laopropytether	<del>                                     </del>	<del> </del>	<del> </del>			110.0					¥10.0	10.0		
1,2-dichloropropane	ļ				<50	<5.0	<5.0				<50	<5.0	<5.0	
1,3-dichloropropana					<50	<5.0	<50				<5.0	<5.0	<50	
2.2 dichloropropara					<5.0	<5.0	<5.0				<5.0	< 5.0	<50	
1.1-dichloroprop-sne	ļ	<b></b>			<5.0	<5.0	<5.0				<50	<5.0	<5.0	
1,3-dichloropropene cis-1,3-dichloropropene		ļ			<10 0 <5.6	<10 0 <5 0	<10.0 <5.0				<10.0	<10.0	<10.0	
tai: 1.3-dichloropropene	<del> </del>				(50	<50	<50 <59				<5.0 <5.0	<5.0 <5.0	<5.0 <5.0	
Ethytonzene	i	<del> </del> -			450	<50	<5.0				<50	<5.0	<5.0	
Heptane														
2-hexanone		<u> </u>			40	<5.0	450				<5.0	< 4.0	<50	
Hexachicrobutadiene	ļ	ļ			0.01>	<10.0	<10.0				<10.0	<10.0	<10.0	
lodomathane laopropyluenzane	<del> </del>		<b> </b>		<10.0	<100 <100	<10.0		·		<10.0	<10.0	<10.0	
- isopropytoluene		<del> </del>	<del> </del>		<10.0	<10.0	<10.0				<10.0 <10.0	<10.0 <10.0	<10.0 <10.0	
Mett-ylene chloxido	<b> </b>		<b>†</b>		₹50	<50	450				<50	<5.0	350	
Methyl ethyl ketone					<10.0	<100	<100				<100	<10.0	<10.0	
MTBE														
Methyl lenbutyl ketone		ļ	ļ										1	
4-metryl-2-pentanone Tetrahydrokaran		<b> </b>	ļ		<50 <50	<5.0	<5.0 <5.0			<u> </u>	<5.0	<5.0	<5.0	
Naphihalene	<del> </del>	<b> </b>			*50	13.0	\$2.0		<del></del>		<50	<5.0	150	
n propytienzene	l	ì			<100	<10.0	<10.0		<b></b>	·····	<10.0	<10.0	<100	
Styrene					<50	<50	<50		<del></del>		<50	<50	<5.0	
1.1,1,2-letrachioroethane	ļ				<50	<5.0	<5.0				<50	<50	<50	-
1,1,2,2-letrachloroethane	ļ				<50	<5.0	<5.0				<50	<5.0	<50	
Tetractilorcathene Toluene	«5 U	<del></del>	<50	<5.0	<50	<5.0	<50		ļ	ļ	<b>450</b>	<5.0	<5.0	
1,2,3-trichlorobenzene	*>0	<5.0		< 5.0	<50 <50	<5.0 <5.0	<5.0 <5.0	<b>₹</b> 5 0	<5.0	<5.0	<5.0 <5.0	<5.0	<80	<5.
1,2,4 trichlorobenzene					450	750	<5.0 <5.0	<del></del>			<5.0 <5.0	<5.0 <5.0	<5.0 <5.0	
1,1,1-trichloroethane	450	<50	<5.0	<50	<5.0	<50	<b>45.0</b>	<5 U	<5.0	<5.0		<5.0	<b>45.0</b>	<5.
1,1,2-trichloroethane					<50	<50	< 5.0				<50	<50	₹5.0	
Trichloroethene					<b>450</b>	<5.0	<5.0				<5 V	<5.0	- 5.0	
Trichkroftyoromethane					<10.0	<10.0	<10.0	<b></b>			<100	<10.0	<10.0	
1.2,3-trichloropropane 1,2,4-transtrytenzene					<50 <100	<5 G	<5.0				<50	< 5.0	<50	
1,3,5 t/methybenzene					<100	<10.0	<10.0				<10.0 <10.0	<10.0 <10.0	<10.0	
Vmyl acetale			<del></del>		<100	₹10.0	410.0				<10.0	410.0	<10.0 <10.0	
Etryl acotate										***************************************		<del>                                     </del>	<u>``'''</u>	
Vrnyl chloride					<10.0	<10.0	4100			·	<10.0	<:0.0	<10.0	
2-chiloroethyl /myl ether												L		, <del></del>
Ethyl methacrylate									,					
o-sylera			ļ		<100	<10.0	<100			·	<100			
ra and p sylene					<50	<5.0	<5.0				< 5.0	<5.0	<5.0	
Kyliones, total					·<50	<5.0	<5.0		<u> </u>	<u> </u>	<5.0	46.0	<5.0	7

Compound	SPRING \$/23/00	SPRING 9/29/00	SPRING 11/35/00	RRUS 1/2/60	RRUS 6/23/08	RRUS 9/29/56	RRUS	RRD8 3/2/00	RRDS 6/23/00	RRD3 9/29/00	RRDS 9/29/00	RRDS 11/JOKO
Acetone												
Acrolein												
Acry/cnitrile		!									<5.0	-5.0
Benzane	<5.0	<5.0	<5.0	<5.0	<5.0	<5.0	<5.0	<5.0	<5.0		15.0	<5.0
Bromotenzene Bromochioromethane												
Bromodichioromethane												
Bromoform												
Bromomethane												
n-buly/benzeno												
sec-buly/benzene												
tert-buty/bonzene												
Carbon disu <b>Ede</b>												
Carbon tetrachloride												
Chlorobenzene												
Chkrod2romomethane		ļ										<u> </u>
Chicroethane												
Chioroform												
Chloromethane		<b> </b>		ļ							<b></b>	
o-chlorotoluene		ļ	<b></b> _		ļ	-						
p-chlorotoluene		<b></b>		<b> </b>			ļi	<b> </b>			ļ	<del> </del>
1,2-dibromo-3-chioropropane	<del> </del>	<b> </b>	ļ	<b> </b> -				<b> </b>			<b> </b> -	<del></del>
1,2-dibromoethane Dibromomethane	<b> </b>	<b>_</b>	<del></del> -	<b> </b>	<b> </b>			<b> </b>			<del> </del>	
1.2-Schlorobenzene	<del> </del>	<del>                                     </del>	<u> </u>	$\vdash$	<del></del>			<b> </b>			<b></b>	<b></b>
1.3-dichlorobenzene	<del> </del>	<del> </del>	<b></b>	<del>  </del>		L		<b></b>			<del> </del>	<del> </del>
p-dichlorobenzene	<del> </del>	1	<del></del> -	t	<del></del> -	ļ		<b></b>			<b></b>	
Dichicrofluoromethane	<del> </del>		<del></del>		<del> </del>						l	·
Dichkrodifluoromi@hane	<del></del>	<del>                                     </del>			<del></del>						<del> </del>	h
1,2-dichloroethane	<del>                                     </del>	<b> </b>			<del> </del>			<b></b>			1	l
1,1-dichloroethane	<b>45.0</b>	<5.0	<50	<5.0	<5.0	<5.0	<5.0	<5.0	<5.0		<5.0	<5.0
1,1-dichloroethene	<5.0		<5.0			<5.0	<50	<5.0	45.0		<5.0	<5.0
cis-1,2-dichloroethers												
trans-1,2-dichloroettiene												
trans-1,4-dichloro-2-butene												
di-isopropylether												
1,2-dichloropropara												
1,3-dichloropropane	<u> </u>		<u> </u>		<u> </u>		<u> </u>					
2,2-dichloropropane	<u> </u>		ļ	<u> </u>	ļ		ļ	ļ			ļ	
1,1-dichloropropene	<b> </b>	<u> </u>	ļ			<u> </u>				ļ	<b></b>	ļ
1,3-dichloropropene				<b> </b>		ļ		<u> </u>	<b></b>		<b></b>	
cis-1,3 dichloropropene		<del> </del> -		<del> </del> -	ļ	<b></b> _			<b> </b>	<b></b>	ļ	<del> </del>
trans-1,3-dichloropropene	·	<del> </del>	ļ	<del> </del>	<b> </b>			<b> </b>		<b> </b> -		
Etrytenzene	ļ	ļ	ļ	ļ	<del> </del>	<u> </u>	ļ	<del> </del>		<b> </b>	<del> </del>	
Heptane 2-hezanone	<b></b>	·	ļ	ļ		<b> </b>	<del> </del>	<del> </del>		<del> </del>		<del> </del>
Hexachlorobutations	<del> </del>	<del> </del>		<del> </del>	<del> </del>	<del> </del>		<del> </del>		<del> </del>	<del> </del>	<del></del>
Iodomethane		<del> </del>		<del> </del>	<del> </del>	<u> </u>		<del> </del>			<del> </del>	<del> </del>
(sopropy/tienzene	<del> </del>	<del> </del>		<del> </del>				<del> </del>			<del> </del>	<del> </del>
p-isopropytichene	<del> </del>	<del> </del>		<del> </del>				·			<del> </del>	<del> </del>
Methylene chloride	1			<b></b>		<del></del>		1		1	1	<u> </u>
Methyl ethyl hatone	L	I	L	L	L			L			I	
MIBE	L										L	
Alethyl isobutyl tretone	<b></b>				<u> </u>		<u> </u>	ļ				
4-methyl-2-pentanone	ļ	ļ	ļ	ļ	ļ	<u> </u>	ļ	<u> </u>	ļ	<b></b>	<b></b>	ļ
Terrahydi ofuran	ļ	ļ	ļ	<b></b>	<b> </b>	<b> </b>		<b>!</b>	ļ	ļ	<del>  </del>	<del> </del>
Naphithalene	<b>}</b>	ļ	<b> </b>	<del> </del>	ļ	<del> </del>	<b> </b>	<b> </b>	<del> </del>		<del> </del>	<del> </del> -
n-propytenzene	<del> </del> -	<del> </del>		}	<del> </del>	<del> </del>	ļ	<del> </del>	<b> </b>	<del></del>	<del> </del>	ļ
Styrene	<del> </del>	<del> </del>	<del> </del>	<del> </del>	<del> </del>	<del> </del>	<del> </del>	<del> </del>		<del> </del>	<del> </del>	<del> </del>
1,1,1,2-let achkroethane	<del> </del> -		<del> </del>	<del> </del>	<del> </del>	<del> </del>	<del> </del>	<del> </del>	<del> </del>	<del> </del>		<del> </del>
1,1,2,2-tet achioroetisane Tet achioroetiene	<del> </del>	<del> </del>	<del> </del>	<del> </del>	<del> </del>	<del> </del>	<del> </del> -	<del> </del>	<del> </del>	<del> </del>	<del> </del>	<del> </del>
Totuene	<5.0	<5.0	<50	450	<5.0	<5.0	<5.0	<5.0	<5.0	<del> </del>	₹5€	<5.0
1.2,3-trichlorobenzene	<del>                                     </del>	<del></del>	<del>                                     </del>	<del> </del>	<del>}</del>	<del></del>	<del>                                     </del>	<del> </del>	1	<del>}</del>	<del> </del>	1
1,2,4-tichlorgbenzene	<del> </del>	<del>                                     </del>	<del> </del>	1	<del> </del>	<del>                                     </del>	<del> </del>	1	<del> </del>	<del> </del>	<del> </del>	·
1,1,1-trichloroethane	<50	<5.0	<5.0	<5.0	<50	<5.0	<5.0	<5.0	₹5.0	<del>                                     </del>	<50	<50
1,1,2-bichloroethane	l	† <del></del>	1	<u> </u>	1	1	T	1	1	1	<del> </del>	†
Inchloroethene	1	1	1	<del> </del>	1	1	1	1	1	1	<del>                                     </del>	1
Trichlorofluorornethane	T	T	I	T	I	Ι	T	1.	1	<del> </del>	7	T
1,2.3 trichlorupropane	I			I	I	I			I	1	]	1
1,2,4-trimethy benzene												L
1,3,5-trimetrytenzeue						L					$\Gamma$	
Vinyl acetata	I		L					L			Ι	<u> </u>
Ethyl acetate	I			L		I		I				T
Vinyl chloride						L		L		L		
2-chkroethylvinyl ether	I		L		J		L	I	L		<u> </u>	T
Ethyl methacrylate			I					L				]
					1							
o xylene												
m and p xylene Xylanes, total							<u> </u>					

Parameter	MW1-S	MW1-S-F	MW1-S	MW1-S	MW1-S-F	MW1-S	MW1-S	MW1-S	MW1-S	HW1-S	MW1-S	MW1-S	MW1-S	MW1-S
	5/7/98	5/7/98	6/17/98	7/24/98	7/24/54	10/26/98	2/16/99	7/26/99	219/59	11/1/99	3/2/00	6/23/00	9/29/00	11/30/00
Aluminum				< 0.10	<0.10	<0.10		< 0.04	<0.04	<0.04	<0.04	<0.04	<0.04	<0.04
Antimony				<0.10	<0.10	<0.10		< 0.001	< 0.001	< 0.001	<0.001	<0.001	0.001	<0.001
Arsenic	<0.0050	< 0.0050		<0.080	40.080	<0.080	<0.080	<0.001	< 0.001	<0.001	<0.001	<0.001	0.001	<0.001
Barium	0.067	0.07		0.064	0.067	0.056	0.069	0.072	0.067	0.066	0.075	0.052	0.069	0.135
Bery/fium				< 0.010	⊴0.010	<0.010	<0.010	<0.001	< 0.001	< 0.001	<0.001	<0.001	< 0.001	<0.001
Boron				<0.1	<0.050	<0.050	₹0.050	< 0.05	0.05	<0.05	0.078	<0.05	<0.05	0.048
Cadrium	<0.0%	<0.010		< 0.020	<0.020	<0.020	<0.020	< 0.001	<0.001	< 0.001	<0.001	<0.001	<0.001	<0.001
Calcium	95	100		100	120	91	100	117	98.3	103	112	100	106	104
Chromium	<0.040	<0.040		<0.020	<0.020	<0.020	<0.020	0.001	<0.001	0.001	0.001	0.001	0.001	6.003
Cotalt				<0.020	<0.020	< 0.020	<0.020	0.001	0.601	0.001	0.004	< 0.001	0.001	0.001
Copper				<0.020	< 0.020	<0.020	< 0.020	0.002	0.002	0.002	0.005	< 0.001	0.001	0.004
iron	<0.050	<0.050	<0.10	<0.10	<0.10	<0.10	<0.10	0.258	0.298	0.153	0.200	0.201	0.220	1.01
Lead	<0.0050	<0.0050		<0.10	<0.10	<0.10	<0.10	<0.001	< 0.001	<0.001	<0.001	<0.001	<0.001	<0.001
Magnesium				48	56	45	49	59.7	48.0	50.7	52.3	49.5	51.5	46.5
Manganese	0.031	0.031	0.029	0.029	0.046	0.025	0.03	0.045	0.039	0.044	0.028	< 0.001	0.032	0.963
Mercury				< 0.0002	< 0.0002	<0.0002	< 0.0002	< 0.0002	<0.0002	<0.0002	< 0.0002	<0.0002	<0.0002	<0.0002
Nickel				<0.058	<0.050	<0.050	<0.050	0.005	0.005	0.006	0.003	0.004	0.005	0.019
Potassium	<1.0	1.1	<20	<20	5.5	<2.0	<2.0	1.04	.89	1.81	1.80	1,13	2.20	<0.05
Selenium	< 0.0050	< 0.0050		<0.15	<0.15	<0.15	<0.15	0.004	0.003	0.003	< 0.001	0.002	<0.001	0.012
Silver			.,	<0.020	<0.020	<0.020	<0.020	<0.001	<0.001	<0.001				
Sodium		12	11	11	12	10	11	12.5	11.3	11.8	12.€	14.5	12.5	11.1
Thalium				<1.0	<1.0	<1.0		<0.001	< 0.001	< 0.001	< 0.001	< 0.001	0.005	<0.001
Vanadium				<0.050	<0.050	< 0.050		0.001	<0.001	<0.001	< 0.001	0.002	0.002	0.008
Zinc				0.08	0.078	<0.020	0.039	0.034	0.032	0.037	0.019	0.007	0.036	0.067
Alkalinity, phenol.				<5.0	<\$.0	<5.0					4.0	<1.0	<1.0	<1.0
Alkalinity, total	254	248		240	260	250		188				240.0	230	246
Alkalinity, bicarbonate	256	248		240	250	250		170		230		227.0	226	238
Ammonia nitrogen				<0.20		<0.20	<0.20	<0.10	0.27	<0.10		40.10	0.23	<0.10
BOD, 5-day				<5.0	<5.0	<5.0		1.0						
Chloride	31	31	: 4		38	30	31	29.6	27.5	27.9		29.2	29.9	28.1
COD				<5.0	<5.0	<5.0	<5.0		<20	<20		<20	<20	<50
Cyanide				<0.00.50	<0.0050	< 0.0050	< 0.0050	<.10	<0.10	<0.10				
Fluoride				<0.20	<0.20	< 0.21	<0.20	0.18	0.15	0.16	0.13	0.17	0.17	0.18
16trate as N	<1.0		6.9		<0.020	9.7	7.5		6.89		3.54	5.49	6.42	5.91
Oil and Grease				<1.5	<1.5	3.1	<5.0							
pH (standard pH units)	7.41		7.3	7.1	7.3	7.4	7.4	7.8	8.2	7.7	7.2	7.5	7.3	8.1
Sulfate	'50	240	170		190	150	180	169	178	178	179	171	165	169
Phenois		<0.020		<0.020	<0.020	<0.020	<0.020			<u> </u>				
Phosphorus, Total		L	<0.10											
TOC				3.9		1	5	6.09	2.32	1.45	3.16	2.73	0.57	0.61
Total Dissolved Solids	574	597	550		550	580	F50	428	440	435	435	422	420	422
Turbidity (nephalometric turbidity units)	1	0.2	<1				1		1			1.37	0.99	

Parameter	MW2-S	MW2-S-F	MW2-S	MW2-S	MW2-S	MW2-S	MWZ-S	MW2-S	MW2-S	MW2-S	MWZ-S	MW2-5	MW2-S	MW3-5
	5/7/98	517798	6/17/98	7/24/98	10/25/98	2/16/99	7/26/99	9/9/99	11/1/99	3/7/00	6/23/09	9/29/00	11/30/00	5/7/98
Vuminum							< 0.04	<0.04	<0.04	< 0.04	<0.04	0.051	0.019	
Antimony							< 0.001	<0.001	< 0.001	< 0.001	<0.001	₹0.001	<0.001	
Arseric	<0.0050	<0.0050					0.003	0.003	0.003	0.002	<0.001	0.002	0.002	0.0176
Barium	0.161	0.158		0.145	0.153		0.172	0.153	0.138	0.165	0,143	0.164	0.118	0.064
Bery <b>sum</b>		•					<0.001	< 0.001	<0.001	< 0.001	<0.001	<0.001	<0.001	
Boron							<0.05	<0.05	<0.05	9.064	<0.05	<0.05	0.015	
Cadmium	<0.005	<0.010					<0.001	< 0.001	<0.001	<0.001	<0.001	<0.001	< 0.001	<0.00
Calcium	81	79		63	86		96.3	88.3	85.6	93.9	86.4	91	82.6	0.1
Chromium	<0.040	<0.040					0.001	< 0.001	0.001	<0.001	0.001	0.002	< 0.003	<0.04
Cobalt							0.003	0.003	0.003	0.002	0.CO1	0.003	0.002	
Copper							< 0.001	0.002	<0.001	0.003	0.004	0.008	< 0.001	
iron	0.198	0.251	0.49	0.2	<0.10		0,568	0.529	0.436	0.371	0.371	0.783	0.190	1.44
Lead	< 0.9050	< 0.0050					< 0.001	0.001	< 0.001	<0.001	0.602	<0.001	<0.001	<0.005
Magnesium							46.4	41.7	40.9	42.0	40.4	41.7	37.6	
Manganese	0,131	0.131	0.062	0.048	0.057		0.055	0.049	0.056	0.050	<0.001	0.088	0,062	0.051
Mercury							Ф.0002	< 0.0002	< 0.0002	< 0.0002	<0.0002	<0.0002	<0.0002	
Nickel							0.004	0.005	0.004	0.003	0.008	0.005	0.003	
Potassium	1.4	1.7	<2.0	<2.0	2.5	<2.0	1.47	1.20	2.80	2.13	6.50	0.920	<0.05	1,400
Selenium	<0.0050	<0.0050					< 0.004	<0.001	<0.001	0.002	<0.001	0.002	0.002	< 0.005
Silver							< 0.001	< 0.001	<0.001					
Sodium	4.2	4.6	4.8	4.3	5.3	4.7	5.15	4.87	4.52	5.30	5.91	5.20	3.40	67
Thallum							<0.001	< 0.001	<0.001	<0.001	<0.001	<0.001	< 0.001	
Vanadium							0.002	<0.001	0.001	<0.001	0.003	0.003	<0.001	
Znc							0.015	0.017	0.014	0.014	0.047	0.100	0.014	
Alkalinity, phenol.				<5.0	<5.0	l				<1.0	<1.0	<1.0	<1.0	
Alkatinety, total	344	344		350	350	i —	281	268	326	339	335.0	332	344	686
Alkalinity, bicarbonate	344	344		350	350		255	255	318	328	331.0	330	323	24
Ammonia nitrogen							< 0.10	0.14	<0.10	< 0.10	<0.10	<0.12	<0.10	
BOO, 5-day			·		<del></del>		1.1	12.7	1.2					
Chloride	<5	<5	<5.0	<5.0	7.7	<5	24.0	1.89	2.16	2.73	2.71	2.92	2.65	13
CO0	<del>-</del>	<u> </u>				<del> </del>	<20	<20	<20	<20	205	77.1	<20	
Cyanide						<del>                                     </del>	<.10							
Fluoride		· · · · · ·	<del> </del>	·	<b></b>	1	0.23	0.10		0.10	0.12	0.12	0.14	
Nétrade as N	6.1	<del> </del>	<1.0	<1.0	<1.0	<del> </del>	<0.10						<0.10	<1.
Oil and Grease	<del></del>		<u> </u>	<del> </del>	<del></del>	1	<.10					l		
pH (standard pH units)	7.21	<del> </del>	7.2	7.1	7.2	7.5	<del></del>			7.0	7.3	7.1	7.8	71 1016
Sulfate	37								34.0	34.4	40.5	40.5	39.4	CA
Phenola			<del></del>	<u></u>	<del></del>				T	ļ —		<del></del>		
Phosphorus, Total		<del> </del>	<0.10	<del> </del>	<u> </u>	<del>                                     </del>	1	1	1	1	<u> </u>			
TOC		<del> </del>	<u> </u>			<del></del>	6.73	1.15	1.05	2.73	3.19	0.63	0.56	
Total Dissolved Solids	395	407	380	370	380	370				337	334	334	332	2,96
Turbidity (nepholometric turbidity s		2.5			<1.0		† <del>*:*</del>	†~~~ <u>**</u>	† <u></u>	<del> </del>	3.49			

Parameter	MW3-S-F	MW3-S	MW3-S	MW3-S-F	MW3-S	MW3-S	MW3-S	MW3-S	MW3.S	MW3-S	WM3-8	MW3-S	MW3-S
	5/7/98	6/17/98	7/24/98	7/24/98	10/26/98	2/16/99	7/26/99	9/9/99	11/1/99	3/2/00	6/23/00	9/29/00	11/30/00
Aluminum							0,29	0.39	0,837	0.406	0.071	0.233	0.028
Anámony							<0.001	<0.001	< 0.001	<0.001	<0.001	<0.001	0.001
Arsenic	0.0304				0.0304		0.131	9.079	0.091	0.065	0.060	0.097	0.175
Barium	0.059		0.049		0.036		0.054	0.046	0.051	0.042	0.034	G.042	0.089
Beryllium							<0.001	< 0.001	<0.001	< 0.001	<0.001	< 0.001	<0.001
Boron							0.05	0.15	0.052	0.184	<0.05	0.052	0.204
Cadmium	< 0.010				<0.010		<0.001	< 0.001	< 0.001	<0.001	< 0.001	< 0.001	0.001
Calcium	8.5		4.4		3		4.44	3.71	5.34	2.40	3.53	6.10	1.30
Chromium	< 0.040				< 0.040		0.003	< 0.001	0.003	0.001	0.003	0.003	0,018
Cobalt							< 0.001	0.002	0.003	< 0.001	<0.001	< 0.001	<0.001
Copper							0.011	0.013	0.014	0.009	0.008	0.045	0.022
iron	< 0.050	<1.0	1		<0.10		0.484	0.399	0.861	0,180	0.152	0.291	0.409
Lead							< 0.001	0.002	0.002	0.001	< 0.001	<0.001	0.001
Magnesium							4.07	2.80	3.43	1.10	2.91	3.50	0.60
Manganese	0.028	< 0.100	0.073		0.031		0.101	0.173	0.328	9,106	< 0.001	0.058	0.041
Mercury							< 0.0002	< 0.0002	< 0.0002	<0.0002	< 0.0002	< 0.0002	0,0015
Nickel							0,014	0.013	0,015	0.008	0.008	0.013	0.038
Potassium	1,200	1,600	1,600	2,000	1,900	2,900	2314	2010	2170	2120	1980	218	2120
Selenium	<0.0050		.,,,,,,,,	7			0.008	a¥ 905	0.007	0.004	0.004	0.005	0.016
Silver							< 0.001	<0.001	<0.001				
Sodium	96	100	98	120	120	120	145	139	134	130	130	138	129
Thatium							< 0.001	< 0.001	<0.001	< 0.001	< 0.001	0.002	<0.001
Vanadium	<b></b>						0.015	< 0.001	0.011	0.003	0.010	0.015	0.029
Zinc	<del> </del>						0.008	0.012	0.014	0.012	< 0.001	0.014	0.015
Alkalinity, phenol.			640		650					662	1165.0	710.0	746
Alkalinity, total	650		1,400		1,500		1590	1600	1472	1540	1507.0	1630.0	1650
Alkalinity, bicarbonate	350		120		200		1450	250	800	293	328.0	807.0	774
Ammonia nitrogen							<0.10	1.60	<0.10	<0.10	< 0.10	0.61	<0.10
BOD, 5-day		1					<1.0	2.9	4.2				
Chloride	418	130	160	170	140	170	181	163	151	160	231	202	203
COD	1						<20	35	<20	36.4	242	68.8	55.9
Cyrnide							<.10	<0.10	<0.10				
Fluoride							3.97	4.00	4.44	3,47	3.11	3.02	3.20
Nitrate as N	1	<1.0	<1.0		<1.0		<0.10	<0.10	0.16	<0.10	<0.10	<0.10	<0.10
Oil and Grease	1						<0.10	<0.10	<0.10				
pH (standard pH units)	T	10.4	10.5	-10.4	4.5. 10.4	10,4	10.9	FEC 113	分/2/10.5	19.4	\$40.010.8	25.1 10.4	10.7
Sulfate	520	1,200	1,400	1100	980	1100	980	1054	1009	931	1153	1170	1242
Phenois	T					<u> </u>	l	]		}			
Phosphorus, Total	T	1.6				I							
TOC	1	T	l		i	T	12.9	22.7	17.0	27.2	17.9	9.26	14.3
Total Dissolved Solids	2,130	3,700	3,900	4,300	4,200	4,300	3776	3462	4350	3980	3150	4400	4390
Turbidity (nephelometric turbidity units)	75	1.57	1		1.8		<u> </u>	1	·		8.10	22,10	<u> </u>

Parameter	MW3-SF	MW4-S	MW4-S-F	NW45	MW4-S	MW4-S	MW4-S	MW4-S	MW4-S	MW4-8	MW4-5	MH4-S	MW4-S G/29/00	MW4-S 11/36/60
	11/30/00	5/7/98	5/7/98	6/17/98	7/24/98	10/26/98	2/16/99	7/26/99	3/9/99	11/1/99	3/2/00	5/23/00		0.017
Aluminum	0.176							0.87	0.96	0.631	0.065	< 0.04	<0.001	<0.001
Antimony	<0.001							<0.001	<0.001	<0.001	<0.001	<0.001	0.001	<0.001
Arsenic	0.106	< 0.0050	< 0.0050					0.003	0.007	0.003	0.001	<0.01	0.107	0,098
Barium	0.070	0.099	0.094		0.084	0.096		0.105	0.094	0.089	0.105	0.068		<0.001
Beryffum	< 0.001							< 0.001	0.004	< 0.001	< 0.001	<0.001	<0.001	0.895
Boron	0.135							1.27	1.10	1.14	1.03	0.08	1.30	
Cadmium	< 0.001	< 0.005	<0.010					<0.001	70.001	<0.001	<0.001	< 0.001	<0.001	<0.001 150
Calcium	0.90	150	160		150	150		177	183	155	168	137	0.002	0.001
Chromium	( 707	<0.040	< 0.040		,			0.003	<0.001	0.002	0.002	0.001		0.001
Cobat	< 0.001			T				0.004	0.005	0.002	0.002	< 0.001	0.001	
Copper	0.018							0.007	0,010	0.005	0.007	<0.001	0.000	0.002
iron	0.366	0.845	< 0.050	<0.10	<0.10	<0.10		2.09	5.39	1.77	0.879	0.323	1.15	0.648
Lead	<0.001	< 0.0050	< 0.0050			11		0.007	0.033	9,006	< 0.001	<0.001	0.001	<0.001
Magnesium	0.40							99.9	8.08	58.1	54.1	52.7	59.3	<b>⊕</b> 54.2
Manganese	0.055	0.359	0.357	0.518	0.670	0.665		0,832	0.952	0.746	0.758	0.407	0.846	0.810
Mercury	<0.0002		-					<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	0.0002	<0.0002
Mickel	0.021							0.010	0.012	0.006	0.005		0.007	0.005
Potassium	2160	250	250	230	200	270	330	234	234	251	214	180	238	233
Selenium	0.012	<0.0050	<0.0050					0.001	0.002	0.002	0.003	<0.001	0.003	0.002
Silver	0.072	-0.0000						<0.001	< 0.001	<0.001				
Sodium	133	25	23	22	20	24	28	27.3	30.3	27.5	24.8	21.0	25.4	24,0
Thalfum	<0.001	<del></del>	<del></del>					< 0.001	< 0.001	<0.001	<0.001		0.001	<0.001
Vanadium	0.016	<del> </del>	<del> </del>	<del> </del>			<b></b>	0.004	<0.001	0.002	0.002		0.003	
Zoc	0.005		<del> </del>					0.037	0.070	0.039	0.021		0,060	
Alkalinity, phenol.	0.000	<del> </del>		<del> </del>	<5.0	<5.0				1	<1.0		<1.0	
AlkaEnity, total	<del></del>	332	328	<del> </del>	360			356	320	384	352	322.0	376.0	
Alkalinity, bicarbonate	<del> </del>	332			360	390		340	310	320	332	320.0	350.0	
Ammonia nitrogen	<del> </del>		<del>                                     </del>	<del> </del>	<del></del>			<0.10	0.23	<0.10	<0.10	<0.10	<0.10	<0.10
BOD, 5-day	<del> </del>	<del> </del>	<del> </del>	<del> </del>		<del></del>		4.3	4.0	<1.0				<u> </u>
Chloride	<del> </del>	110	115	84	83	90	120	81.0	92.6	99.2	\$6.1		84,3	
COD	<del> </del>	<del>                                     </del>	<del> </del>	<del>                                     </del>			<del>                                     </del>	7	₹20	<20	<20	52.6	57.1	54.4
	<del> </del>	<del> </del>	<b> </b>	<del>                                     </del>	<del> </del>	<del> </del>	1	<0.10	<0.10	Ø).10			ſ	
Cyanide	<del> </del>	<del> </del> -	<del> </del>	<del>                                     </del>			<del> </del>	0.21	0.22	0.22	0.16	0.25	0.25	0.24
Fluoride	<del> </del>	<1.0		<1.0	<1.0	<1.0	·	<0.1	0.11	<0.10	<0.10	<0.10	<0.10	< 0.10
Nitrate as N	+	1	<del> </del>	1-1.0	†	1	1	<0.1				1		
Oil and Grease	- <del> </del>	7.16		7.3	7.2	7.3	3 7.			7.	7.	7.4	7.3	
pH (standard pH units)	<del> </del> -	53/2			510					502	455	275	443	41
Sulfate		1- 337	1	+	† <del></del>	<del> </del>	1	1	1	1	1	T	T	T
Phenols		<del> </del>	<del> </del>	<0.10	<del> </del>	<del>                                     </del>	1	<del> </del>	<b>†</b>	1	T	T	1	T
Phosphorus, Total	<del> </del>	<del> </del>	<del> </del>	<del></del>	<del> </del>	<del> </del>	<del> </del>	11.5	3.3	4.5	8.31	9,01	2.4	0 2.5
TOC	<del> </del>	1.420	1,410	1,300	1,200	1,40	1,70			·			88	2 88
Total Dissolved Solids Turbidity (nephelometric turbidity units)		1,42%	<0.			<1.		<del>\                                    </del>	1	<del> </del>	1	4.2		

Parameter	MW4-SF	MW5-S	MW5-S	MW5-S	MW5-S	MW5-S	MWSS-F	MW5-S	MW5-S	WW5-SF	MNV6-S	FAW6-S	MV-6-S	MW6-S
	11/30/00	7/26/99	9/9/99	11/1/99	3/2/60	6/23/00	6/23/00	9/29/00	11/30/00	11/30/00	7/26/99	6/9/99	11/1/99	3/2/00
Aluminum	<0.04	0.17	0,17	0.161	0.286	0.041	<0.04	0.151	0.195	<0.04	3.01	0.08	0.116	0.166
Antimony	<0.001	< 0.001	<0.001	<0.001	<0.001	<0.001	<0.001	< 0.001	<0.001	<0.001	<0.001	< 0.001	<0.001	< 0.001
Arsenic	< 0.001	0.001	<0.001	0.003	0.005	<0.001	< 0.001	0.001	< 0.001	<0.001	0.004	<0.001	0.001	0.003
Barium	0.095	0.199	0.069	0.095	0,307	0.078	0.086	0.131	0.087	.082	0,350	0.331	0,271	0.276
Be:y@um	< 0.001	<0.001	< 0.001	<0.001	< 0.001	< 0.001	< 0.001	<0.001	< 0.001	< 0.001	<0.001	< 0.001	<0.001	0.001
Boron	0.979	0.52	0.55	0.527	0.700	<0.05	<0.05	0.572	0.328	0.300	0.51	0.63	0.688	0.431
Cadmium	<0.001	<0.001	<0.001	<0.001	0.002	<0.001	< 0.001	<0.001	<0.001	<0.001	< 0.001	< 0.001	<0.001	0.001
Calcium	143	274	291	279	278	204	218	177	175	172	345	24.6	232	398
Chromium	0.002	0.004	0.001	0.003	0.006	0.002	0,003	0.004	0.002	0.001	0.007	< 0.001	0.002	0.005
Cobalt	0.001	< 0.001	<0.001	< 0.001	0.008	<0.001	< 0.001	< 0.001	< 0.001	<0.001	0.020	0.008	0.009	0.030
Copper	0.001	0.004	0.008	0.007	0.017	<0.001	<0.001	0.003	0.004	0.004	0.024	0.006	0.004	0.022
Iron	0.398	0.913	0,816	0.718	2.24	0.351	0.381	0.804	0.644	0.354	5.34	0.693	1.06	5.65
Lead	< 0.001	0.001	0.001	<0.001	0.001	< 0.001	< 0.001	<0.001	0.001	<0.001	0.015	< 0.001	<0.001	<0.001
Magnesium	53.6	89.0	34.6	101	202	28.9	58.3	58.7	62.1	52.7	78.0	31.4	51.9	
Manganese	0.780	0.081	0.064	0.047	1.25	< 0.001	<0.001	0.044	0.117	0.083	3.11	0.694	2.44	2.13
Mercury	< 0.0002	< 0.0002	< 0.0002	<0.0002	<0.0002	<0.0002	<0.0002	< 0.0002	< 0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002
Nickel	0.006	0.005	0.004	0.003	0.021	0.001	0.002	0.002	0.002	0.002	0.033	0.017	0.017	0.038
Potassium	222	23.5	61.9	59.5	341	23.9	221	74.5	17.7	145	84.6	70.2	68.6	59.8
Selenium	0.004	0.003	0.001	0.004	0.006	<0.001	<0.001	0.001	0.001	< 0.001	0.003	0.001	0.002	0.005
Silver		< 0.001	< 0.001	< 0.001							< 0.001	<0.001	< 0.001	
Sodium	22.6	22.8	78.5	55.3	34.0	47.2	33.0	18.6	9.	17.1	35.1	55.0	27.3	33.4
Thailium	<0.001	< 0.001	<0.001	<0.001	<0.001	< 0.001	< 0.001	< 0.001	< 0.001	<0.001	< 0.001	<0,001	< 0.001	0.001
Vanadium	0.002	0.007	<0.001	0.005	0.023	0.005	0.007	0.007	0.003	0.002	0.023	< 0.001	0.003	0.902
Zinc	0.020	0.011	0.020	0.012	0.047	< 0.001	< 0.001	0.009	0.016	0.009	0.041	0.010	0 011	0.063
Alkalinity, phenol.	T				<1.0	<1.0		<1.0	<1.0					<1.0
Alkalinity, total		324	372	586	386	432.0		490.0	502		369	354	390	
Alkalinity, bicarbonate		314	350	540	377	427.9		375.0	479		360	325	350	379
Ammonia nitrogen	1	-0.10	<0.10	<0.10	<0.10	<0.10		0.19	<0.10		<0.10	<0.10	<0.10	<0.10
800, 5-day	T	4.6	1.6	<1.0							5.3			
Chloride	T	21.6	111	95.5	125	36.9		25.2	23.9		187	142	105	
COD		<20	26	<20	<20	57.9		25.7	203		<20			
Cyanide	1	<0.10	<0.10	<0.10							<0.10		<0.10	
Fluoride	T	0.13	0.06	0.09	0.06	0.13		0.13	0.12		0.29	0.17	0.22	0.19
Nitrate as N		1.31	6.24	2.00	<0.10	<0.10		<0.10	0.10		0.10			
OI and Greas		<0.10	<0.10	<0.10							<0.10			
pH (standard pH units)		7.6	7.5	7.6	6.9	7.1	7.3	7.3	7.7		7.4	7.9	7.8	7.8
Sulfate	T	220	571	578	650	278		198	203		214	332	314	242
Phenois	T	T	T	T	T									
Phosphorus, Total	1		1	T		T	T							
TOC	1	21.6	29.0	17.9	< 0.50	9.86		1.41	1.46		21.1	13.2	5.59	
Total Dissolved Solids	]	668	1090	1016	1215			633	845		867	885	820	789
Turbicity (nephelometric turbidity units)	1	Ī		T	1	3.62		180.00						

Parameter	MW6-S	MW6-S	MW6-S	MAY8-SF	HW7-S	MW7-S	MW7-S	MW7-8	HW7-S	MW7-S	MW7-S	MW8-S	MW8-S	WWW.S
	6/23/00	9/29/00	11/30/00	11/30/00	7/26/99	9/9/99	11/1/99	3/2/00	6/23/00	9/29/00	11/30/00	7/26/99	9/9/99	11(1/99
Numinum	<0.04	0.127	1.74	0.101	.83	6.30	0.443	0.662	<0.04	0,292	0.090	0.51	0.29	0.665
Antimony	<0.001	<0.001	< 0.001	<0.001	< 0.001	<0.001	< 0.001	0.003	<0.001	< 0.001	<0.001	<0.001	<0.001	< 0.001
Arsenic	<0.001	0.001	0.003	6.003	0.028	0.025	0.044	0.017	0.007	0.013	0.026	0.001	< 0.001	0.001
Barium	0.143	0.251	0,181	0.120	0.045	0.047	0.046	0.055	0.033	0.044	0.025	0.073	0.666	0.056
Beryllium	< 0.501	<0.001	<0.001	< 0.001	<0.001	<0.001	<0.001	< 0.001	< 0.001	< 0.001	<0.001	< 0.001	< 0.001	<0.001
Boron	<0.05	0.292	0.364	0.437	0.25	0.19	0.216	1.95	<0.05	0.204	9.195	0.25	0.25	0.225
Cadmism	<0.001	< 0.001	<0.001	<0.001	<0.001	<0.001	<0.001	< 0.001	<0.001	< 0.001	<0.001	< 0.001	<0.001	<0.001
Calcium	149	233	258	235	23.9	14.7	11.9	43.4	40.8	29.0	14.2	265	226	223
Chromium	0.002	0.603	0.003	0.002	0.004	0.002	0.002	0.005	0.002	0.001	0.002	0.004	0.003	0.006
Cobalt	<0.001	0.007	0.011	0.068	0.001	< 0.001	<0.001	0.001	<0.001	< 0.001	<0.001	0.002	0.002	0.002
Copper	0.001	0.012	0,020	0.005	0.005	0.007	8.008	0.003	<0.001	0.006	0.004	0.015	0.014	0.011
Iron	0.284	0.770	4.68	0.759	1,98	9.789	0,629	1.32	0.109	0,683	0.142	1,60	1,10	1.57
Lead	<0.001	<0.001	0.008	<0.001	0,003	0,002	0.002	<0.001	< 0.001	0.001	<0.001	0.003	0.004	9.603
Magnesium	20.3	47.7	57.6	47.3	10.8	5.80	4.55	13.6	15.1	10.6	4.60	110	94.4	94.3
Mangariese	0.094	1,81	1.79	1.67	0.154	0.133	0.085	0.157	0.002	0.103	0.024	0.363	0.233	0.255
Mercury	<0.0002	< 0.0002	<0.0002	<0.0002	< 0.0002	< 0.0002	<0.0002	< 0.0002	<0.0002	<0.0002	<0.0002	< 0.0002	<0.0002	<0.0002
Nickel	0.003	0.009	0.023	0.012	0.012	0.014	0,014	0.007	0.002	0.011	0.009	0.010	0.010	0.008
Potassium	31.2	49.5	39.7	43.1	656	717	793	503	406	823	727	265	263	265
Selenium	<0.001	0.003	0.004	0.003	0.001	0,002	0.002	0.003	<0.001	0.003	0.002	0.003	0.003	0.003
Sitver	<0.001				<0.001	<0.001	<0.001		L			<0.001	<0.001	<0.001
Sodium	53.6	64.9	55.6	57.8	37.6	42.9	49.9	37.4	25.8	44.2	48.8	8.2	38.6	40,1
Thallium	<0.001	< 0.001	<0.001	< 0.001	< 0.001	<0.001	<0.001	<0.001	< 0.001	<0.001	<0.001	<0.001	<0.001	<0.001
Vanadium	0.005	0.006	0.013	0.001	0.016	0.612	0.022	0.015		0.013	0.006	0.003	<0.001	0.003
Zinc	< 0.001	0.013	0.023	0.004	0.027	0.015	0.016	0.027	<0.001	0.023	<0.001	0.044	0.027	0.042
Alkalinity, phenol.	<0.001	<1.0	<1.0					41	<1.0	92.0				
Alkalinity, total	417.0	384.0	420		396	466	534	348	341.0	384.0	486	271	270	314
Alkalinity, bicarbonate	410.0	350,0	357		285	265	475	154	333.0	254.0	297	265	265	306
Ammonia nitrogen	<0.10	<0.10	<0.10		<0.10	<0.10			<0.10	0.94	<0.10			
BCO, 5-day		<u> </u>			3.1					<b></b>		3,1		<1.0
Chloride	71.1	246	188		70.1	90.2	97.7	56.3	48.8	105	110	119	121	113
COD	100	123	94.1	<u> </u>	<20	<20			10.5	34,3	<20			
Cyanide		L			<0.10	<0.10			L	<u> </u>		<0.10		
Fluoride	0.41	0.27	0.30	<u> </u>	1.84	1.95	2.43	0.80	0.65	9,30	1.84	0.26	0.16	
Nitrate as N	<0.10	<0.10	<0.10	<u> </u>	<0.10				<0.10	<0.10	<0.10			
Oil and Grease		<u> </u>		<u> </u>	<0.10				ļ	<u></u>	ļ	<0.10		
pH (standard pH units)	7.1				10.0				8.9					
Sulfate	106	186	254	<u> </u>	401	475	494	345	336	148	474	618	818	763
Phenois		L		<u> </u>		<u> </u>	<u> </u>	<u> </u>		<u> </u>	ļ		<u> </u>	<b></b>
Phosphorus, Total	1					<u> </u>	<u> </u>	<u> </u>		<u> </u>		<u> </u>	ļ	
TOC	13.2	2.01			17.2	10.7	15.2		18.0	7.00		9,80		
Total Dissolved Solids	566	842	830	L	1047	1294	1423	801	802	1156		1135	1100	1191
Turbidity (nephelometric turbidity u	nits) 2.69	200.00	. k			1	1	I	0.6	23.00	4	l .	1	i

Parameter	MW8-S	MW8-S	MWS-S	MYV8-S	MWB-SF	MW9-S	MW9-S	MW9-S	MW9-S	MW9-S	MW9-S	KW9-S
	3/2/00	6/23/00	9/29/00	11/30/00	11/30/00	7/26/99	9/9/99	11/1/99	3/2/90	6/23/00	9/29/00	11/20/00
Aluminum	0.435	< 0.04	0.109	0.233	< 0.04	0.52	0.02	0.389	0.698	<0.04	0.149	0.019
Antimony	<0.001	< 0.001	<0.001	<0.001	< 0.001	0.003	0.002	< 0.001	0.002	< 0.001	<0.001	<0.001
Arsenic	6.005	< 0.001	0.001	< 0.001	< 0.001	0.013	0.005	800.0	0.009	0.004	0.003	0.002
Bacture	0.172	0.044	0.055	0.057	0.058	0.086	0.065	0.060	0.058	0.047	0.041	0.033
Bery/Sum	0,002	< 0.001	< 0.001	<0.001	<0.001	<0.001	< 0.001	<0.001	<0.001	<0.001	< 0.001	<0.001
Boron	0.253	<0.05	0.199	0.232	0.211	< 0.05	0.15	0.067	0.070	< 0.05	0.032	0.075
Cadmium	0.003	<0.001	<0.001	<0.001	< 0.001	< 0.001	< 0.001	< 0.001	<0.001	<0.001	<0.001	<0.001
Calcium	268	224	234	222	210	175	118	394	207	108	110	97.4
Chromium	0.040	0.007	0.005	0.006	0.004	0.004	< 0.001	0.002	0.007	0.002	0.002	<0.003
Cobalt	0.035	< 0.001	<0.001	<0.001	<0.001	0.004	0.002	0.003	0.004	< 0.001	0.601	<0.001
Copper	0.113	< 0.001	0.007	0.097	0.004	0.019	0.012	0.034	0.036	0.003	0.013	0.010
kon	5.30	0.415	0.549	9.953	0.210	2.00	0,582	1.29	5.10	0.231	0.827	0.13
Lead	0.004	< 9.001	<0.001	0.002	< 0.001	0.008	0.002	0.007	< 0.001	<0.001	<0.001	<0.001
Magnesium	125	92.4	93,3	91.1	85.4	86.8	72.2	81.2	104	72.3	73.5	56,2
Manganese	4.30	< 0.001	0.024	0.187	0.065	0.472	0.202	0.212	0.224	0.040	0.192	0.101
Mercury	< 0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	0.0003	<0.0002	<0.0002	< 0.0002	< 0.0002	<0.0002
Nickel	0.058	0.002	0.003	6.005	0.004	0.020	0.018	0.010	0.012	0.003	0.007	0.00
Potassium	217	174	254	280	274	364	378	394	477	428	445	42
Sejenjum	0.014	0.001	0.002	0.006	0.003	0.002	0.001	0.004	0.009	0.003	0.004	0.005
Sitver	1					< 0.001	<0.001	<0.001				
Sodium	43.1	43.4	52.4	51,4	50.7	89.2	88.0	91,9	103	101	07.6	93.7
Thallium	0.002	<0.001	0.005	0.001	0.001	<0.001	<0.001	< 0.001	<0.001	0.003	<0.001	<0.00
Vanadium	< 0.001	0.004	0.003	0.003	0.001	0.004	<0.001	0.002	0.002	0.005	0.004	0.00
Zinc	0.363	< 0.001	0.025	0.021	0.018	0.029	0.014	0.025	0.050	< 0.001	0.015	0.00
Alkalinity, phenol.	<1.0	<1.0	<1.0	<1.0				1	<1.0	<1.0	<1,0	<1.
Alkalinity, total	284	316.0	274.0	278		465	398	236	334	459.0	458.0	476
Alkalinity, bicarbonate	271	314.0	262.9	257		391	355	220	287	423.0	375.0	454
Ammonia nitrogen	<0.10	<0.10	<0.10	<0.10		0.30	0.51	<0.10	<0.10	<0.10	0.53	<0.1
BOO, 5-day						2.9	4.4	1.2				
Chloride	162	181	179	198		76.9	90.1	109	129	152	108	110
COD	30.3	158	<20	<20		<20	<20	<20	<20	216	<20	<21
Cyanide						<0.10	<0.10	<0.10				
Fluoride	0.16	0.26	0.22	0.21		0.31	0.23	0.26			0.31	0.34
Nitrate as N	1.68	2.95	3.01	1.86		0.10	<0.10	<0.10	0.14	0.12	<0.10	<0.1
Oil and Grease						<0.10	<0.10	<0.10				
pH (standard pH units)	7.1	7.2				7.9	8.2				7.4	7.
Sulfate	832	686	847	865		575	640	668	622	703	717	68
Phenois							<u> </u>					
Phosphorus, Total				L			L					
TOC	14.4	15.1	2.05			7.72	4.10		8.46	9.90	1.90	2.2
Total Dissolved Solids	1161	1125	1200	1252		1064	1172	1257	1365	1333	1250	127
Turbidity (nephelometric turbidity units)	1	3.03	21.00	1	1	1	I	}	1	2.26	18.20	

Parameter	MW1-D	MW1-D-F	MW1-D	MW1-D	MW1-D-F	Kw1-D	MW1-D	MW1-D	MW1-D	MW1-D	MW1-5	MW1-D	MW1-D
	5/7/98	5/7/98	8/17/98	7/24/98	7/24/95	10/26/98	2/16/99	7/26/99	9/9/99	11/1/98	3/2/00	6/23/00	9/29/00
Aluminum			<0.04	<0.10	<0.10	< 0.10		<0.04	0.08	0.109	0.111	<0.04	0.049
Antimony			<0.601	<0.10	<0.10	<0.10		< 0.001	<0.001	<0.001	<0.001	<0.001	<0.001
Arsenic	<0.0050	<0.0050	<0.001	<0.080	<0.080	<0.080	<0.080	<0.001	<0.001	<0.001	0.001	< 0.001	0.001
Barium	0.063	0.049	< 0.001	0.044	0.051	0.647	0.046	0.047	0.053	0.047	0.055	0.030	0.048
Beryllium			< 0.001	<0.010	< 0.010	<0.010	<0.010	<0.001	< 0.001	<0.001	< 0.001	< 0.001	<0.001
Boron			<0.05	<0.1	< 0.050	< 0.050	< 0.050	<0.05	<0.05	<0.05	0.051	<0.05	<0.05
Cadmium	<0.005	<0.010	<0.001	<0.020	< 0.020	<0.020	<0.020	<0.001	< 0.001	<0.001	<0.001	<0.001	<0.001
Calcium	93	71	<0.05	80	100	86	79	92.3	87.6	84.9	89.4	51.8	82.4
Chromium	< 0.040	< 0.040	<0.003	< 0.020	<0.020	<0.020	<0.020	.002	<0.001	0.002	0.001	<0.001	0.002
Cobalt			< 0.001	<0.020		< 0.020	<0.020	<0.001	0.004	0.003	0.007	<0.001	<0.001
Copper			< 0.001	<0.020	< 0.020	< 0.020	<0.020	0.002	0.003	0.004	0.003	<0.001	0.007
Iron	0,872	< 0.050	<0.005	<0.10	<0.10	<0.10	<0.10	.273	0.568	0.519	1.18	0.132	0.214
Lead	<0.0050	< 0.0050	<0.001	<0.10	<0.10	<0.10	<0.010	<0.001	0.001	<0.001	0.002	< 0.001	<0.001
Magnesium			< 0.005	39	46	44	41	49.7	45.5	44.0	121	36.2	43.7
Manganese	0.071	<0.010	< 0.001	<0.010	<0.010	<0.010	<0.010	0.007	0.050	0.053	0.219	< 0.001	0.002
Mercury			<0.0002	< 0.0002	<0.0002	<0.0002	< 0.0002	< 0.0002	<0.0002	<0.0002	< 0.0002	< 0.0002	< 0.0002
Nickel			<0.001	<0.050	<0.050	<0.050	<0.050	0.002	0.003	0.003	0.003	<0.001	0.002
Potassium	1.5	1,4	<0.05	<2.0	8,5	<2.0	<2.0	1.57	1.34	1.89	2.90	1.20	1.90
Selenium	<0.0050	<0.0050	<0.001	<0.15	<0.15	< 0.15	< 0.15	< 0.004	0.001	0.001	0.004	0.001	0.008
Silver				<0.020	<0.020	<0.020	<0.020	< 0.001	< 0.001	< 0.001			
Sodium	3.1	3.2	<0.001	3.1	3.4	4.2	3.3	3.65	3.57	3.90	3.80	3.00	3.60
Thatium			<0.001	<1.0		<1,0		<0.001	< 0.001	<0.001	0.002	<0.001	0.003
Vanadium			< 0.001	< 0.050	< 0.050	<0.050	< 0.020	0.002	< 0.001	0.002	0.001	0.002	0.004
Zinc			< 0.001	<0.020	<0.020	<0.020	<1.0	0.011	0.013	0.018	0.017	<0.001	0.028
Alkalinity, phenol.			<1.0	<5.0	<5.0	<5.0					<1.0	<1.0	<1.0
Alkalinity, total	348	336	<1.0	350	350	350		282	268	320	314	325.0	312.0
Alkalinity, bicarbonate	348	336	<1.0	350	350	350		250	260	310	296	308.0	310.0
Anmonia nitrogen			<0.10	<(1E	S <0.20	<0.20		<0.10	0.11	<0.10	< 0.10	<0.10	0.29
BOD, 5-day				<5.0	<5.0	<5.0		2.5	3.7	2.4			
Chloride	<5	<5	<0.05		<5.0	<5.0	<5.0	2.23	1.91	2.55	2.58	4.86	2.6
COD			<20	<5.0	<5.0	<8.3	₹5.0	<20	<20	<20	<20	52.6	43.7
Cyanide				<0,0050	<0.0050	< 0.0050	< 0.0050	<.10	<0.10	<0.10			
Fluoride			<0.05	<0.20	<0.20	<0.21	<0.20	0.18	0.13	0.14	0.11	0.14	0.13
Nitrate as N	<1.0		<0.10			<1.0	l	0.20	018	0.19	0.14	0.13	0.2
Oil and Grease				<1.5		<2.7	<5.0	<.10	<0.10	<0.10			<u> </u>
pH (standard pH units)	7,31	<del> </del>	<del> </del>	7.1		7.4	7.3	8.1	8.1	7.7	7.2	7.4	7.
Sulfate	37	43	<0.10		39	35	37	37.7	37.0	37.6	35.6	40.7	40.
Phenols	<del>-</del>	<u> </u>		<0.020		<0.020	<0.020				l		1
Phosphorus, Total		<u> </u>		1			1	l				<u> </u>	1
TOC			<0.50	1.5	2.3	<1.0	6.0	4.13	2.79	3.92	4.59	3.10	<0.5
Total Dissolved Solids	396	390	<1.0		390	400	370			326		323	
Turbidity (nephelometric turbidity units)		<0.1		<b></b>	<u> </u>	·	<del></del>			1		0.00	

Parameter	MW1-D	MW2-D	MW2-D-F	MW2-D	MW2-D	MW2-D	MW2-D	MW2-D	MW2-D	IAW2-D	MW2-D	MW2-D	MW2-D
	11/30/00	5/1/98	5/7/98	6/17/94	7/24/98	10/26/98	2/15/99	7/26/99	9/9/99	11/1/09	3/2/00	6/23/00	9/29/00
Aluminum	<0.04							< 0.04	<0.04	<0.04	0.040	< 0.04	0.053
Antmony	<0.001							<0.001	<0.001	< 0.001	<0.001	<0.001	<0.001
Arsenic	<0.001	0.0101	0.0111					0.015	0.011	0.015	0.019	0.005	0.011
Barlum .	0.036	0.131	0.141		0.147	0.162		0.179	0.157	0.160	0.185	0.151	0.160
Beryllium	<0.001							< 0.001	< 0.001	< 0.001	<0.001	<0.001	<0.091
Boron	0.009							< 0.05	<0.05	<0.05	0.059	∢0.05	<0.05
Cadmium	<0.001	<0.005	< 0.010					<0.001	<0.001	<0.001	<0.001	<0.001	<0.001
Calcium	69.1	92	92		100	99		115	97.2	13	111	105	104
Chromium	<0.003	<0.040	<0.040					0.002	< 0.001	0.001	0.001	0.001	0.001
Cobait	<0.001							0.004	0.003	0.003	0.002	0.001	0.001
Copper	0.001							0.001	0.002	0.001	0.005	0.001	0.003
iron	9.217	0.764	0.752	1.1	1.5	1.6		2.50	2.47	2.45	2.17	1.07	2.15
Lead	<0.001	< 0.0050						0.002	<0.001	<0.001	< 0.001	0.001	<0.001
Magnesium	38.5							51.0	42.1	45.2	48.3	45.6	44.9
Manganese	0.001	0,615	0.668	0.545	0.618	0.578		0.287	0.264	0.208	0.199	0.104	0.184
Mercury	< 0.0002							<0.0002	<0.0002	<b>40.0002</b>	<0.0002	<0.0002	<0.0002
Nickel	0.001							0.003	0.003	0.002	0.003	0.005	0.002
Potassium	<0.05	1.4	1.7	<2.0	<2.0	<2.0	<2.0	1.38	0.92	1.40	2.40	1.72	2.30
Selenium	0.004	< 0.0050	< 0.0050					< 0.004	<0.001	< 0.001	< 0.001	< 0.001	<0.001
Silver								< 0.001	< 0.001	<0.001			
Sodium	2.46	12	12	9.7	9.2	9,7	9.8	10.8	9.64	10.1	11.0	11.1	11.1
Thatium	<0.001	- 12						<0.001	< 0.001	< 0.001	< 0.001	< 0.001	< 0.001
Vznadium	0.001							0.003	<0.001	0.001	< 0.001	0.003	0.004
Zinc	0.004							0.032	0.012	0.010	0.013	0.023	0 023
Alkalinur, phenol.	₹1.0				<5.0	<5.0					<1.0	<1.0	<1.0
Alkalinity, total	328	404	404		400	400		316	324	382	344	376.0	366.0
Alkalinity, bicarbonate	315	404	404		400	400		3:0	320	350	329	357.0	355.0
Ammonia nitrogen	<0.10							<0.10	0.10	<0.10	<0.10	<0.10	<0.10
BOD, 5-day								<1.0	6.3	<1.0			
Chloride	2.46	10	19	23	22	22	22	20.1	19.8	18.9	20,3	21.9	31.9
COD	<20							<20	<20	<20	<20	68.4	160
Cyanide								<.10	<0.10	<0.10			
Fluoride	0.17	<del>`</del>		<del> </del>				.21	0.16	0.17	0 12	0.19	0.17
Nitrate as N	0.18			<1.0	<1.0	<1.0		<0.16	<0.10	0.34	<0.10	<0.10	<0.10
Oil and Grease	<del> </del>	<u></u>		·				<.10	<0.10	<0.10			
pH (standard pH units)	8.0	7.15	<del> </del>	7.2	7.2	7.3	7.4	7.6	8.2	7.7	7.2	7.3	7.2
Sulfato	39.9		40		38	29		29.6	25.2	30.6		29.2	30.4
Phenois		<del></del>	<del></del>	1	† <del></del>	<u></u>	·			·	T		
Phosphorus, Total	<del> </del>	<del> </del>	<b> </b>	<0.10	<del> </del>	<del>                                     </del>	<del> </del>		l		<b></b>		
70C	0.28			1	<del> </del>	<del> </del>	<b> </b>	8.71	6.11	1.05	5.15	10.1	<0.50
Total Dissolved Solids	318		472	470	420	450	440	388	398	390	393	1301	396
Turbidity (nephelometric turbidity units)	<del> </del>	<del> </del>	7.0		† <del></del>	<1.0		l	1	<del>                                     </del>	T	3.93	

Parameter	MW2-D	MW2-DF	MW2-DF	MW3-D	MW3-D-F	MW3-D	MW3-D	MW3-D	MW3-D	MW3-D	MW3-D	MW3-D	MW3-D
	11/30/50	11/30/00	2/14/61	5/7/98	5/7/98	6/17/98	7/24/98	10/25/98	2/16/99	7/26/99	\$/9/99	11/1/99	3/2/00
Aluminum	< 0.04	<0.04	< 0.04							<0.04	0.055	<0.04	0.050
Ansmorry	<0.001	<0.001	<0.001							<0.001	<0.001	<0.001	<0.001
Arsenic	0.008	0.005	0.013	0.0080	0.0101					0.019	0.006	0.022	0.017
Barium	0.230	0.140	0.160	0.134	0.133		0,139	0.139		0.163	0,053	0.134	0.159
Berylium	<0.001	< 0.001	< 0.001							40.001	<0.001	<0.001	< 0.001
Boron	0.022	0.011	<0.05							<0.05	<0.05	<0.05	0.076
Cadmium	<0.001	<0.001	< 0.001	< 0.005	<0.10					<0.001	<0.001	<0.001	<0.001
Calcium	100	100	29.4	93	8.6		99	9.5		112	41.8	98.9	104
Chromium	0.004	0.001	=0.001	<0.040	< 0.040					0.002	<0.001	0.002	< 0.001
Cobalt	0.002	0.001	0.001		315.13					0.002	0.002	0.003	0.001
Copper	0.002	0.0010	0.002							0.001	0.002	6.002	0.002
iron	2.55	1.59	2.24	0.919	0.560	0.80	0.90	1.1		1.79	0.830	2.03	1.57
Lead	<0.001	<0.001	<0.001		0.500	0.00	0.50			<0.001	<0.001	<b>40.001</b>	<0.001
Magnesium	42.5	42.4	43.8	0.904	0.846	0.547	0.504	0.477		49.2	17.3	41.7	42.5
Manganese	0.300	0.170	0.136	- 3,17.7	4.010	- 0.011	0.501	V.177		0.314	0.087	0.282	0.172
Mercury	<0.0002	<0.0002	<0.0002							<0.0002	<0.0002	<0.0002	<0.0002
Neckel	0.003	0.002	0.002							0.008	0.001	0.002	0.002
Potassium	1.30	1.80	1.68	1.8	1,9	<2.0	4.1	2.6	2.0	3.58	2.30	3.82	2.90
Selenium	0.007	0.002	<0.003	<0.0050	<0.0050		7.1			< 0.004	<0.001	<0.001	0.001
Silver	1			-0.0000	-0.0050					<0.001	- CO.001	<0.001	0.001
Sodium	9.50	9.60	6,92	8.9	8.4	9.6	9.1	7.4	6.6	6.76	2.90	6.43	
Thallium	<0.001	<0.001	<0.001					- '^	0.0	<0.001	<0.001	<0.001	6.90
Vanadium	0.008	0.001	<0.001							0.003	0.001	0.001	<0.001 <0.001
Zinc	0.013	0.059	0.007							0.009	0.006	0.025	0.009
Alkalinity, phenol.	<1.0					~	<5.0	<5.0		0.003	0.000	0.023	<1.0
Alkalinity, lotal	396			384	316		380	350		310	364	366	350
Alkalinity, bicarbonate	380			384	316		380	340		295	348	345	347
Ammonia nitrogen	<0.10									<0.10	0.14	<0.10	<0.10
BOD, 5-day			<del></del>		··					2.5	5.5		<b>40.10</b>
Chloride	<0.05			7	7	8.5	8.8	8.3	8.6	7.27	7.78	<1.0 7.13	7.81
COD	35,3									<20	137	<del>7.13</del> <20	<20
Cyanide										<10	<0.10	<0.10	٠,٨
Fluoride	0.30									0.28	0.22	9.20	
létrate as N	<0.10	<del>i</del>		<1.0		<1.0	<1.0	<1.0		<0.10	<0.10	<0.10	0.17
Oil and Grease										<.10	<0.10	<0.10	<0.10
pH (standard pH units)	7.9			7.24		7.1	7.1	7.3	7.3	8.1	7.2		
Suttate	30.6			39	37	42	36	28	30	27.5		7.8	7.0
Phenois									30[	- 47.5	30.9	23.6	27.0
Phosphorus, Total	t			<del> </del>		<0.10							
TOC	0.74	<del></del>								<del></del> -	<del>- ,,</del> ,		
Yotal Dissolved Solids	306			432	431	450	420	410	400	10.6 359	0.61	1.56	2.63
Turbidity (nephelometric turbidity units)			<del></del>		4.0	<del>430</del> <1	720	<1.0	100	359	353	357	362

Parameter	MW3-D	MW3-D	MW3-D	MW3-DF	MW4-D	MY4-D-F	MW4-D	MW4-D	MW4-D-F	MW4D	MW4-D	MW4-D	MW4-D
	6/23/00	9/29/00	11/30/00	11/30/00	5/7/98	C/7/98	6/17/98	7/24/98	7,7,4/98	2/16/09	7/26/99	9/5/99	11/1/99
Aluminum	<0.04	10.04	<0.04	<0.04							<0.04	0.06	<0.04
Албтопу	<0.001	<0.001	<0.001	<0.001							<0.001	<0.001	<0.001
Arsenic	0.009	<0.001	0.015	0.016	0.0053	< 0.0050					0.004	0.003	0.004
Barium	0.118	<0.001	0.127	0.120	0.125	0.126		0.099	0.088		0.086	0.081	0,076
Bery'sum	<0.001	<0.001	<0.001	<0.001							<0.001	<0.901	<0.001
Boron	<0.05	<0.05	0.023	0.020							1.10	0.75	0.957
Cadmium	<0.001	< 0.001	<0.001	<0.001	<0.005	<0.010					< 0.001	<0.001	<0.001
Calcium	95,3	< 0.05	99.4	95.8	160	160		200	180		272	177	191
Chromium	0.001	< 0.001	< 0.003	<0.003	<0.040	< 0.040					0.002	< 0.001	0.001
Cobalt	0.001	<0.001	0.002	< 0.001							0.003	0.002	0.002
Copper	<0.001	<0.001	<0.001	<0.001							0.004	0.003	0.004
iron	1,32	< 0.005	1.46	2.84	1.12	1.08	0.64	0.49	0.39		0.957	.02	0.667
Lead	< 0.001	<0.001	<0.001	< 0.001	< 0.0050	< 0.0050					<0.001	<0.001	<0.001
Magnesium	40.1	<0.005	40.8	40.5			***************************************				85.7	21.0	73.4
Manganese	0,119	< 0.001	0.157	0.138	2.74	2.78	1.4	1.1	0.907		1.13	1.16	1.09
Mercury	<0.0002	<0.0002	<0.0002	<0.0002							<0.0002	<0.0002	< 0.0002
Nickel	0.002	< 0.001	0.002	0.003							0.017	0.013	0.018
Potassium	4.92	<0.05	6.00	5.50	180	160	160	150	160	180	214	193	209
Selenium	<0.001	<b>30.001</b>	<0.001	<0.001	<0.0050	<0.0050					< 0.001	0.001	0.001
Silver											< 0.001	< 0.001	<0.001
Sodium	6.12	<0.001	6.10	5.80	27	28	22	21	21	22	24.3	67.4	23.2
Thalium	< 0.001	< 0.001	<0.001	< 0.001							0.001	0.001	<0.001
Vanadium	0.004	< 0.001	< 0.001	<0.001							0.003	< 0.001	0.002
Zinc	<0.001	< 0.001	<0.001	< 0.001	i						0.142	0.079	0.166
Alkalinity, phenol.	<1.0	<1.0	<1.0					<5.0	<5.0				
Alkalinity, total	360.G	<1.0	364	<u> </u>	404	404		390	410		413	400	400
Alkaśnity, bicarbonate	345.0	<1.0	327		404	404		390	410		390	375	370
Ammonia nitrogen	<0.10	<0.10	<0.10								<0.10	<0.10	<0.10
BOD, 5-day											5.3	5.5	<1.0
Chloride	8.56	<0.05	8.42		93	93	71	76	76	88	93.8	101	84.8
COD	232	<20	<20								<20	<20	<20
Cyanide			ſ								<.10	<0.10	<0.10
Fluoride	0.17	<0.05	0.20		1						0.36	0.19	6.24
Nitale as N	<0.10	<0.10	<0.10		<1.0	1	<1.0	<1.0	<1.0		0.10	<0.10	<0.10
Oil and Grease				· · · · · · · · · · · · · · · · · · ·							<0.10	<0.10	
pH (standard pH units)	7.2	<u> </u>	7.9	1	6.93	l	7.0	6.9	7.2	7,1	7.8	8.3	7.4
Sulfate	28.7	<0.10	32.3		520	510	530	550	390	559	514	537	484
Phenois	l	1	1	1		<u> </u>	i						
Phosphorus, Yolal		T	1	1	I	[	0.13	1					
TOC	5.89	<0.50	1.36	1	l	<u> </u>					8.79	11.1	<0.50
Total Dissolved Solids	357	<1.0	357		1,330	1,320	1,300	1,200	1,300	1,400	966	1034	950
Turbidity (nephelometric turbidity units)	3.46	1.40	<b>X</b>	1		10	<1		<1.0		I	1	

Parameter	MW4-D	MW4-D	MW4-D	MW4-D	MWS-D	14W5-D	MWS-D	85₩5-D	MW5-D	MW5-D	MWS-D	MAKE-D	MWE-D
	3/2/00	6/23/00	9/29/00	11/30/00	7/26/99	9/9/99	11/1/99	3/2/00	6/23/00	9/29/00	11/30/00	7/25/99	8/8/99
Aluminum	<0.04	<0.04	0.198	0.012	0.36	0.06	0.050	0.080	<0.04	0.105	0.023	0.11	0.0
Antimony	<0.001	<0.001	<0.001	< 0.001	<0.001	<0.001	<0.001	< 0.001	<0.001	< 0.001	49.001	<0.001	€0.60
Arsenic	0.003	0.001	0.002	0.002	0.002	<b>⊴0.001</b>	0.002	0.003	<0.001	0.002	0.001	0.905	0.00
Barium	0.082	0.066	0.080	0.089	0.121	0.062	0.053	0.087	0.043	0.052	0.050	0.095	0,062
Beryllium	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	< 0.001	C.001	<0.001	< 0.001	<0.001	<0.001	<0.00
Boron	1.17	0.06	1.13	0.960	0.45	0.40	0.350	0.407	<0.05	0.390	0.468	0.56	0.40
Cedmium	<0.001	< 0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	< 0.001	<0.001	<0.001	< 0.001	<0.00
Calcium	205	182	206	186	235	213	214	247	205	224	218	198	167
Chronium	0.002	0.001	0.002	0.001	0.004	<0.001	0.002	8.004	0.001	0.004	0.001	0.003	<0.00
Cobalt	0.002	0.001	0.001	0.002	0.015	0.008	0.008	9.044	0.004	0.307	0.008	0.007	0.00
Copper	0.064	<0.001	0.030	0.003	0.028	0.011	0.007	0.033	<0.001	0.013	0.004	0.037	0.00
Iron	0.618	0.355	1.06	0.386	0.993	0.544	0.385	2.24	0.383	0.496	0.307	1.43	0.41
Lead	< 0.001	< 0.001	0.006	< 0.001	0.002	< 0.001	<0.001	<0.001	<0.001	<0.001	<0.001	0.010	<0.00
Magnesium	77.1	70.8	76.3	72.6	85.5	30.3	76.9	82.7	71.3	79.2	73.1	66.6	56.3
Manganese	1.29	0.533	0.979	0.804	0.475	0.213	0.242	0.70	0.094	0.215	0.284	0.378	0.101
Mercury	<0.0002	< 0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	<0.0002	< 0.0002	<0.0002	< 0.0002	< 0.000
Nickel	9,017	0.012	0.017	0.019	0.013	0.007	0.008	0.020	0.004	0.005	9.006	0.008	9.004
Potassium	248	198	233	204	456	449	416	438	401	359	341	184	173
Selenium	0.006	<0.001	0.002	0.004	0.064	0.002	0.003	0.004	< 0.001	0.004	0.004	0.002	<0.00
Silver					< 0.001	<0.001	<0.001					<0.001	<0.00
Sodium	25.5	23.6	28.2	22.2	22.1	75.0	31.4	33,4	30.3	31.2	28.9	20.7	18.4
Thalfurn	0.001	0.002	0.001	0.003	0.003	0.001	0.002	< 0.001	0.001	< 0.001	0.001	<0.001	< 0.00
Vanadium	0.002	0.005	0.005	0.003	0.005	<0.001	0.002	0.005	0.004	0.004	0.001	0 003	<0.00
Zinc	0.537	0.122	0.218	0.284	0.032	0.015	0.016	0.058	<0.001	0.021	0.004	0.038	0.01
Alkalinity, phenol.	<1.0	<1.0	<1.0	<1.0				<1.0	<1.0	<1.0	<1.0		i
All-atinity, total	400	398.0	378.0	394	339	338	346	408	338.0	334.0	350	303	170
Alkalinity, bicarbonate	385	388.0	365.0	\$77	323		325	385	332.0	320.0	340	290	16
Ammonia nitrogen	<0.10	<0.10	< 0.10	<0.10	<0.10	<0.10	<0.10	< 0.10	<0.10	0.10	<0.10	<0.10	1.0
BOD, 5-day	l				2.3	2.5	1.2					3.0	<1.
Chloride	116	108	107	110	159	148	150	164	164	157	160	92.2	80.6
COD	<20	168	94.3	<20	<20	<20	<20	<20	31.6	114	<20	<20	<20
Cyanide					<0.10	<0.10	< 0.10					< 0.10	<0.10
Fluorida	0.15	0.35	0.27	0.25	0.15	0.09	0.11	0.08	0.13	0.13	0.13	0.18	0.1
Nárate as N	<0.10	<0.10	<0.10	<0.10	1.07	1.11	1.18	0.50	1.09	0.80	1.66	1.63	1.71
Oil and Grease					<0.10	<0.10	<0.10					<0.10	<0.1
pH (standard pH units)	7.2	7.1	7.1	7.5	7.9	5.4	7.8	7.0	7.2	7.4	7.5	7.9	8.
Sulfate	549	516	561	531	871	830	881	850	315	873	834	533	497
Phenois	1					1							<del></del>
Phosphorus, Total	T	i — —				1							i
TOC	7.91	10.8	1.76	2.24	10.8	4.34	4.03	7.76	10.8	1.29	2.19	2 69	2.8
Total Discolved Solids	1032	971	986	926	1376	1383	1384	1360	1322	1270	1271	873	876
Turbidity (nepholometric turbidity units)	<u> </u>	1,20	40.60		1	T			1.37	5.40			·

Parameter	MWS-D	MW6-D	MW6-D	MW6D-F	MNV6-D	MWS-D	MW7-D	MW7-D	MW7-D	MW1-D	MW7-D	MW7D-F	MW7-D
r a siste	11/1/09	3/2/00	6/21/00	6/23/90	9/29/00	11/30/00	7/26/99	879/39	11/1/99	3/2/00	6/23/00	6/23/00	9/29/00
Aluminum	<0.04	0.063	<0.04	4004	0.097	<0.04	<0.04	<0.04	0.045	<0.04	<0.04	<0.04	0.040
Artimony	<0.001	<0.001	<0.001	<0.001	< 0.001	<0.001	< 0.001	< 0.001	< 0.001	<0.001	<0.001	<0.001	<0.601
Arsenic	0.004	0.013	<0.001	<0.001	0.003	0.001	0.004	0.011	0.017	0.012	0.007	0.005	0.008
Barium	0,059	0.078	0.046	0.043	0.058	0.049	0.091	0.160	0.141	0.174	0.033	0.113	0.131
Bery/Eum	< 0.001	<0.001	<0.001	< 0.061	<0.001	<0.001	<0.001	<0.001	<0.001	< 0.001	<0.001	<0.001	<0.001
Boron	0.449	0.540	<0.05	<0.05	0.499	0.394	< 0.05	<0.05	<0.05	<0.05	<0.05	<0.05	<0.05
Cadmium	<0.001	<0.001	< 0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001
Calcium	174	188	180	18¢	214	201	106	90.4	\$3.5	103	93.5	94.3	69.9
Cironium	0.002	C.005	0.002	0.002	0.002	<0.003	0.001	< 0.001	0.002	0.005	03002	0.001	0.001
Cobalt	0.003	0.027	<0.001	<0.001	0.003	0.001	0.004	0.002	0.002	0.002	< 0.001	<0.001	<0.001
Copper	0.003	0.039	< 0.001	<0.001	0.010	0.003	0.009	0.004	0 004	6.002	<0.001	<0.001	0.003
Iron	0.352	4,54	0.344	0.319	0.913	0.183	0.206	2.41	2.19	3.81	1.37	1.68	1.88
Lead	<0.001	<0.001	<0.001	< 0.001	0.001	<0.001	< 0.001	<0.001	<0.001	0.002	<0.001	<0.001	<0.001
Magnesium	59.9	61.8	58.3	60.0	74.0	69.2	47.2	39.0	40.6	421	39.9	40.7	37.7
Manzanese	0.113	0.278	<0.001	<0.001	0.056	0.061	0.600	0.310	0.689	0.290	0.087	0.090	0.140
Mercury	<0.0002	<0.0002	< 0.0002	<0.0002	< 0.0002	<0.0002	<9.0002	<0.0002	<0.0002	<0.0002	<0.005	<0.0002	<0.0002
läckel	0.004	9.011	0.002	0.002	0.004	0.003	0.006	0.005	0.005	0.005	0.001	0.001	0.002
Potassium	152	166	229	235	181	127	2.43	2.21	3.37	1.10	1.81	2,90	3.50
Selenium	0.001	0.006	<0.001	<0.001	0 002	0.001	<0.001	<0.001	< 0.001	0.002	<0.001	<0.001	<0.001
Silver	<0.001						<0.001	<0.001	< 0.001				
So/sium	18.3	19.5	18.8	21.4	21.8	18.3	8.2	6.98	6.57	8.90	5.70	6.00	5.70
Thelium	10.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.901	<0.001	<0.001	<0.001	<0.001	<0.001
Vanadium	9.002	0.005	0.003	0.004	0.004	0.001	0.003	<0.001	0.002	<0.001	0.004	0.001	0.003
Zinç	0.012	0.048	0.013	0.007	0.020	<0.001	0.018	0.028	0.016	0.030	<0.001	< 0.001	0.013
Alkafinity, phenol		<1.0			<1.0	<1.0	<u> </u>			<1.0			<1.0
Alkalinity, total	330	280			304.0	318	326		376	362	384.0		354.6
Alkalinky, bicarbonate	300	266			287.0	297	308		350	345	347.0	<b></b>	342.0
Ammoria ritrogen	<0.10	<0.10	<0.10		<0.10	<0.10			<0.10		<0.10		0.16
BOD, 5-day	<1.0	L		ļ		L	3.3		<1.0				
Chloride	85,6	\$3.8	125	ļ	126	110	3.71		4.72	5.47	5.59	\	5.53 <20
COO	<20	<20	78.9	<u></u>	<20	124	<20		<20		<20	ļ	~20
Cyanide	⊴0.10				<del></del>	<u> </u>	<0.10				ļ	<del> </del>	0.13
Fluoride	0.13				0.16		0.23		0.20	0.16			
Nitrate as N	1.72	1.70	1.73		0.07	0.64					<0.10	<del> </del>	<0.10
Of and Grease	<0.10		<u> </u>	ļ	ļ	ļ	<0.10				<del> </del>	<del> </del>	7.3
pH (standard pH units)	7.4			7.2					7.7				
Sulfate	487	485	517		592	534	20.7	16.6	18.6	20.0	20.2	<b></b>	21.2
Phenois	ļ	<b> </b>	ļ	<b> </b>	<b></b>		<del> </del>	<del> </del>	ļ	<b> </b>	<del> </del>	<del> </del>	ļ
Phosphorus, Total	ļ	<u> </u>	<u> </u>		<del> </del>	ļ	<del> </del>	<del> </del>	<del></del>	320	13.7	<del> </del>	0.00
тос	1.62		11.6	ļ	1,43				4.4			ļ	0.67
Total Dissolved Solids	852	<b>b37</b>	943		939		34(	355	353	347	350		344 11.0
Turbidity (nephelometric & rbidity units)	<u> </u>	<u> </u>	4.8	i	11 00	1	<u> </u>	.L	L	ــــــــــــــــــــــــــــــــــــــ	1 10.	1	11.0

Parameter	MWT-D	MY7-DF	MAN-D	MWI-D	MW8-D	MW3-D	MW4-D	HW8D-F	INVS-D	MW8-D	MWII-DF	MW9-D	MW9-D
į	11/39/99	11/39/0	1/1/1/19	9999	11/1/99	3/2/00	6/23/00	6/23/00	9/29/00	11/33/00	11/30/00	7/25/29	9/2/7
Uurninuss	0.004	0.033	4004	<0.04	< 0.04	< 0.04	40.74	<0.04	0.087	<0.04	<0.04	<0.04	<0.0
Interiory	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	400 11	<0.001	<0.001	<0.001	<0.001	<0.001	<0.0
Arsenic	0.011	0.010	0.016	0.017	0.024	0.019	i. )08	0.010	0.013	0.011	0.007	<0.001	0.0
Parkura Parkura	9.131	150	0.175	0.245	6.230	0.265	0.1.)5	0.214	6.253	6.238	0.215	0.140	0.18
Beryllium	40.001	6.01	40.001	<0.001	<0.001	40.001	40.001	<0.001	40.001	<0.001	<2.001	<0.001	<0.0
Beron	0 017	0.026	40.05	<0.05	<0.05	40.0%	<0.05	<0.05	<0.05	0.030	0.27	<0.05	<0.
Cadmium	<0.001	<0.001	<0.001	₹0.001	€.001	₹0.001	<0.001	45.001	<0.001	<0.001	< 0.001	<0.001	9.0
Calcium	90.2	88.5	<b>∮</b> 23	921	31.8	95.8	91.8	90.0	97.3	\$2.2	90.5	108	92
Chromium	0.001	0.005	0.002	0.008	0.001	6	0.001	0.002	4.001	0.004	<0.001	0.001	<0.0
Cobatt	<0.001	40.001	0.004	0.004	0.003	0.002	<0.001	40.001	<0.001	0.001	0.001	0.003	0.0
Copper	0.002	(,001	0.015	0.005	6.003	0.004	<0.001	<0.001	0.003	0.003	<0.001	0.002	00
roa	1.80	1.72	1.13	3.70	3,46	3.45	1.30	2.27	2.37	1.63	1.37	0.239	5.4
Lead	40 001	40.001	0 001	0.003	Ø.001	<0.001	<0.001	<0.001	40.001	<b>49.001</b>	<0.001	<0.001	40.0
Magnesium	37.7	36.2	45.9	39.3	4C.1	40.4	33.2	38.1	42.3	38.0	36.7	46.8	4
Vanganese	0.152	6.346	1.31	0.613	1.21	0.602	0.292	0.222	0.281	0.296	0.287	1.40	0.2
Mercury	40.0002	1,0002	40,0002	0.0010	40.0002	<0.0002	<0.0002	<0.0002	40.0002	<0.0002	<0.0002	<0.0002	40.00
lictal	0.002	6.0vi	0.000	0.012	0.065	0.005	0.002	0.001	8.004	0.000	0.005	0.010	8.0
Potassium	4.00	400	212	2.12	1.83	2.30	3.00	4.54	2.05	1.10	<0.05	7.69	4.
Selerium	0.003	0.007	<0.001	40 501	<0.001	0.004	<b>49 001</b>	<0.001	0.002	0.001	0.002	<0.001	<00
Siver		T	40.001	<0.001	40.001				<0.001			<0.001	<0.0
Sodium	4.90	5.10	9.54	9,15	7.47	7.22	6.70	6.20	6.80	6.00	5.49	7.16	6.
Thatium	<0.001	4001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	< 0.001	<0.001	9
Vanadum	0.062	0.003	0 002	40 001	0 00%	40 001	0.901	0.005	9.004	0.002	0.002	0.003	<0.
Zinc	0.005	0.008	0.016	0.012	0.021	0.015	<0.001	<0.001	0.014	<0.001	0.005	0.011	0.
Allafrity, pherol.	<1.0					6	<1.0		<1.0	<1.0			
Allralinity, total	362	<del> </del>	341	362	380	364	376.0	1	374.0	378		310	
Alkafinity, bicarbonate	341		323	335	360	341	359,1		354.0	259	i	295	
Ammonia nitrogen	<0.10	1	<0.10	<0.10	<0.10	4.10	<010		<0.10	<b>40.10</b>		<0.10	0
900, 5-day			4.9	2.2	1.8	-						3.6	
Chloride	5.52		1.04	1.67	1.90	1.56	1.73		1.57	1.70		3.26	3
COO	47.1	1	<20	<20	<20	<b>(20</b> )	163		<20	<20		<20	
Cyanics		l	<0.19	<0.10	<0.10	1						<0.10	4
Fluoride	0.20	<del> </del>	0.27	0.21	0.22	0.19	0.24		0.22	0.25		0.23	
Nitrata as 11	Ø.10		<0.10	<0.10	<0.10	4010	<0.10		<0.10	<0.10		0.33	4
Oil and Grease		<u> </u>	<0.10	<310	<0.10		<b></b>	<del></del>				<0.10	4
pH (standard plif units)	80		8.1	8.2	7.7		7.3		7.2	7.7		7.9	
Sulfate	20 3		9.15	9.5/	10.7	10.6	8.81		9.79	9.83		12.4	,
Prenois		<del> </del>	1	1	1	1	<del> </del>		<u> </u>	1	<u> </u>	l	
Phosphorus, Yotal			T	<del>                                     </del>	1	1	<b>†</b>	T		<del> </del>	1		·
TOC	0.66	l	10.0	11.1	<030	2.31	15.3	<b>1</b>	0.63	0.83	1	7.90	,
Total Dasytred Solids	344		347						339			357	
Turbidity (neohelor, retric turbidity units)		T	1	1	†	†****	14.7	·	19.00		<del> </del>		<del> </del>

Parameter	MW9-D	G-WWG	MW9-D	MA-D	MW9-D	MW9-DF
T an all the	11/1/99	3/2/06	6/23/00	2/29/00	11/10/00	11/30/00
Aurinum	1.77	0.510	<0.04	1.20	0.045	0.004
Албитопу	<0.001	<0.001	<0.001	<0.001	40.001	<0.001
Arsenic	0.020	0.008	0.005	0.015	0.908	0.006
Barium	0.215	0.188	0.136	0.225	0.157	0.161
Beryllium	<0.001	< 0.001	<0.001	0.001	< 0.001	<0.001
300	<0.05	< 0.05	< 0.05	< 0.05	0.021	0.019
Cadmium	<0.001	<0.001	< 0.001	<0.001	<0.001	<0.001
Calcium	3.10	102	94.0	113	92.3	93.5
Chromium	0.013	6.006	0.002	0.006	<0.003	<0.003
Cobatt	0.003	0.001	<0.001	0.01	< 0.001	<0.001
Copper	0.004	0.002	<0.001	0.003	0.002	0.001
Irca	10.3	5.16	2.09	8.93	2.23	2.12
Lead	0.006	< 0.001	< 0.001	0.013	0.002	<0.001
Magnesium	42.6	42.8	39.7	43.7	38.7	38.7
Manganese	0.574	0.554	0.057	0.225	0.124	0.127
Mercury	<0.0002	<0.0002	< 0.0002	<0.0002	< 0.0002	<0.0002
Mickel	0.016	0.006	0.001	0.008	0.001	0.001
Potassium	3.10	1.50	3.10	3.40	3.60	2.80
Selanium	9.001	0.003	< 0.001	0.002	0.003	0.001
Sit/er	<0.001			< 0.001		
Sodem	6.87	7.10	6.63	7.50	6.10	5.90
Thatsum	<0.001	< 0.001	< 0.001	<0.001	<0.001	<0.001
Vanadium	0.5.77	0.003	0.004	0.007	0.001	0.001
Zinc	0.123	0.046	<0.001	0.076	0.008	800.0
Alkalinity, phenol.		<1.0	<1.0	<1.0	<1.0	
Alkalinity, total	370	352	371.0	378.0	378	
Alkalinity, bicarbonate	350	339	365.0	332.0	367	
Ammonia nkrogen	<0.10	<0.10	<0.10	0.30	<0.10	
BOO, 5-day	3.6					
Chloride	3.22	4.29	4.92	4.62	4.25	
COD	<20	<20	516	<20	23.5	
Cyaride	<0.10		۰			
Fluoride	9.20	0.17	0.18	0.19	0.21	
Notate as N	<0.10	<0.10	<0.10	<0.10	<0.10	
Oil and G.: AS6	<0.10					
pH (standard pH units)	7.9		7.2	7.8		
Sulfate	15.5	17.5	15.2	25.0	15.9	
Phenois				L		L
Phosphorus, Total						
ICC	7.09	8.54	12.1	0.82		<u> </u>
Total Dissolved Solids	356	348	347		347	L
Turbidity (nephelometric turbidity units)			3.8	3		

Parameter	MWI-WT	MW1-WT	MANUS MIT	LOWE WET	24194 197	MWI-WT	MWI-WT	MWI-WTF	1000	Leun um	2411/2 11/2	MW2-WT	14302 1125
1 at anieczi	7/26/99	9/9/99	11/1/99	MW1-WT 3/2/60	6/23/00	9/29/00	11/30/00	11/30/00	MW2-WT 7/26/99	MW2-WT 9/9/99	MW2-WT 11/1/99	3/2/00	6/23/00
Aluminum	0.27	<0,04	< 0.04	< 0.04	0.080	0.226	0.074	< 0.04	0.31	0.16	9.275	0.312	<0.04
Antimony	< 0.001	<0.001	< 0.001	< 0.001	< 0.001	< 0.001	< 0.001	<0.001	< 0.001	< 0.001	< 0.001	< 0.001	< 0.001
Arsenic	0.003	0.002	0.001	0.002	0.002	0.001	< 0.001	< 0.001	0.002	0.002	0.002	0.001	< 0.001
Barium	0.122	0.155	0.140	0.183	0.250	0.202	0.142	0.079	0.106	0.087	0.067	0.070	0.049
Beryllium	< 0.001	< 0.001	< 0.001	< 0.001	< 0.001	< 0.001	< 0.001	<0.001	<0.001	< 0.001	< 0.001	< 0.001	<0.001
Boron	< 0.05	<0.05	0.061	0.078	< 0.05	< 0.05	0.031	0.016	< 0.05	< 0.05	< 0.05	0.080	< 0.05
Cadmium	< 0.001	< 0.001	<0.001	< 0.001	< 0.001	< 0.001	< 0.001	< 0.001	< 0.001	<0.001	< 0.001	< 0.001	< 0.001
Calcium	97.7	107	99.1	116	96.4	129	125	120	328	282	238	415	174
Chromium	0.002	< 0.001	0.001	0.001	0.001	0.601	< 0.003	< 0.003	0.004	< 0.001	0.002	0.003	0.002
Cobalt	0.007	0.008	0.001	0.004	0.004	0.005	0.005	0.002	0.002	0.002	0.002	0.005	< 0.001
Copper	0.012	0.011	0.005	0.005	0.001	0.008	0.004	< 0.001	0.002	0.012	0.002	0.003	0.001
iron	0.730	<del></del>	0.175	0.966	0.304	0.620	0.582	0.048	1.40	1.09	0.844	1.55	0.346
Lead	0.003	0.004	<0.001	< 0.001	0.001	0.001	< 0.001	< 0.001	0.006	0.004	0.002	0.002	< 0.001
Magnesium	18.2	32.7	35.7	39.6	38.8	49.4	46.8	45.8	121	107	91.3	121	64.4
Manganese	2,04	2,64	0,912	2.23	1.88	1.78	2.36	1.25	0,223	0.181	0.286	0.089	< 0.001
Mercury	< 0.0002	<0.0002	< 0.0002	< 0.0002	< 0.0002	< 0.0002	< 0.0002	< 0.0002	< 0.0002	< 0.0002	< 0.0002	< 0.0002	< 0.0002
Nickel	0.023	0.016	0.002	0.010	0.022	0.018	0.017	0.002	0.002	0.002	0.002	0.002	
Potassium	2.43	3.06	3.63	1.90	3.30	4.60	3.30	3.90	1370	1480	1010		0.003 488
Selenium	< 0.004	< 0.001	< 0.001	< 0.001	< 0.001	0.005	0.002					816	
Silver	<0.001	<0.001	< 0.001	< 0.001	₹0.001	ψ.υσ	0.002	< 0.001	0.005	0.008	0.008	0.005	0.002
Sodium	64.9	38.5	35.1	34.4	26.2	31.5	26.1	26.9	< 0.001	< 0.001 62.4	<0.001 48.7	44.3	31.4
Thallium	< 0.001	<0.001	<0.001	0.002	< 0.001	0.001	< 0.001						
Vanadium	0.005	< 0.001	0.002	0.002	0.007	0.007	0.004	<0.001 <0.001	< 0.001	0.001	< 0.001 0.002	< 0.001 0.004	< 0.001 0.004
Zinc	0.010	0.011	0.008	0.002	< 0.001	0.007	0.006	0.001	0.003	0.012			
Alkalinity, phenol.	1		1 0.000	0.012	<1.0	<1.0	<1.0	0.001	0.013	0.012	0.012	0.017	< 0.001
Alkalinity, total	201	346	322	215	270.0	290.0	280		384	362	398	<1.0 308	<1.0
Alkalinity, bicarbonate	198	330	320	188	265.0	276.0	272		365	355	390		478.0
Ammonia nitrogen	<0.10	< 0.10	< 0.10	< 0.10	< 0.10	0.14	<0.10		< 0.10	< 0.10	< 0.10	298	471.0
BOD, 5-day	2.2	35.9	1.8	9.4	V 0.10	0.14	20.10					<0.10	<0.10
Chloride	1.55	22.1	17.9	18.8	18.2	19.6	18.6		3.1 297	6.6 678	<1.0 464	383	777
COD	<20	44	<20	25.6	<20	42.9	<20						131
Cyanide	<.10	<0.10	< 0.10	< 0.10		42.9	<u> </u>		<20	<20	<20	< 20	353
Fluoride	0.13	1.13	0.12	0.10	0.13	0.13	0.15		<.10 0.23	< 0.10 0.09	< 0.10	0.00	
Nitrate as N	<0.10	<0.10	0.12	<0.10	< 0.10	<0.10	<0.10		0.23	0.85	0.11	0.06	0.24
Oil and Grease	<.10	<0.10	< 0.12	0.121		70.10	₹0.10		<.10		1.59	1.40	0.20
pH (standard pH units)	7.5	8.3	7.8	8.0	7.4	7.1	7.6		7.6	< 0.10	< 0.10		
Sulfate	80.6	86.1	95.7	120	175	238	253		£121	8.1 1619	7.7 1300	7.2	7.2
Phenois									1101	1019	1300	1150	771
Phosphorus, Total													
тос	8.06	22.3	4.04	7.44	10.4	1.28	1.43		10.9	12.7	14.0	17.0	
Total Dissolved Solids	320	430	407	390	420	485	471		3168	3412	2910	17.9 1968	6.65 1407
Turbidity (nephelometric turbidity									3100	7714	4710	1708	144/
units)	<u> </u>	L			8.50	3.00							2.54

							* 41X/2 11/T	1411/2 1177	MW3-WT	MW3-WT	MW3-WTF	MW4-WT
Parameter	MW2-WT 9/23/00	MW2-WT 11/30/00	MW2-WTF 11/30/00	MW3-WT 7/26/99	MW3-WT 9/9/99	11/1/99	3/2/00	MW3-WT 6/23/00	9/29/00	11/30/00	11/30/60	7/26/99
Aluminum	0.190	0.060	0.060	0.25	0.16	0.091	0,149	0.041	0.069	0.047	0.042	< 0.04
Antimony	< 0.001	< 0.001	< 0.001	0.001	< 0.001	0.001	< 0.001	<0.001	< 0.001	0.003	< 0.001	< 0.001
Arsenic	0.001	0.001	0.001	0.006	0.009	0.008	0.004	0.002	0.002	0.020	0.007	0.003
Barium	0.088	0.114	0.103	0.114	0.045	0.041	0.018	0.032	0.055	0.142	0.048	0.158
Beryllium	< 0.001	< 0.001	< 0.001	< 0.001	< 0.001	< 0.001	< 0.001	< 0.001	< 0.001	< 0.001	.<0.001	< 0.001
Boron	< 0.05	0.056	0.060	0.11	0.12	0.112	0.156	< 0.05	0.066	0.452	0.129	1.57
Cadmium	< 0.001	< 0.001	< 0.001	< 0.001	< 0.001	< 0.001	< 0.001	< 0.001	< 0.001	< 0.001	< 0.001	< 0.001
Calcium	269	299	300	86.7	16.5	11.9	2.20	9.10	18.0	15.2	3.50	440
Chromium	0.002	0.002	0.002	0.007	0.001	0.003	0.002	U.003	0.001	0.014	0.002	0.002
Cobalt	0.002	< 0.001	< 0.001	0.002	0.002	0.001	< 0.001	< 0.001	< 0.001	0.002	< 0.001	0.010
Copper	0.006	0.007	0.006	0.014	0.011	0.011	0.006	0.007	0.099	0.022	0.012	0.003
fron	0.946	0.352	0.263	1.03	0.228	0.131	.220	0.097	0.152	0.407	0.106	1.87
Lead	0.003	< 0.001	< 0.001	0.006	0.002	< 0.001	0.002	0.002	0.001	< 0.001	<0.00i	< 0.001
Magnesium	96.9	109	111	26.9	11.7	7.72	1.40	10.7	17.4	12,4	12.2	71.1
Manganese	0.435	0.031	0.027	0.228	0.151	0.158	0.056	< 0.001	0.091	0.095	0.110	2.20
Mercury	< 0.0002	< 0.0002	< 0.0002	< 0.0002	< 0.0002	< 0.0002	< 0.0002	< 0.0002	< 0.0002	0.0009	< 0.0002	< 0.0002
Nickel	0.010	0.008	0.007	0.010	0.010	0.006	0.002	0.006	0.003	0,026	0.010	0.019
Potassium	669	783	956	892	1210	1080	767	675	560	1050	1090	261
Selenium	0.006	0.006	0.007	0.015	0.006	0.009	0.003	0.005	0.005	0.023	0.005	0.002
Silver	1		1	<0.001	<0.001	< 0.001						< 0.001
Sodium	33.7	51.2	51.0	75.7	85.5	78.8	63.7	81.0	63.6	95.4	94.8	21.3
Thallium	0.002		0.001	< 0.001	< 0.001	< 0.001	0.003	< 0.001	0.005	< 0.001	< 0.001	< 0.001
Vanadium	0.006	<del></del>	0.001	0.014	0.018	0.019	0.009	0.007	0.005	0.064	0.014	0.005
Zinc	0.018	0.011	0.012	0.012	0.007	0.005	0.00	0.015	0.013	0.016	0.006	0.035
Alkalinity, phenol.	<1.0	<1.0					160	<1.0	52.0	300		
Alkalinity, total	376.0	430		507	922	852	530	437.0	368.0	760		363
Alkalinity, bicarbonate	360.0	422		433	245	790	16	428.0	355.0	420		350
Ammonia nitrogen	< 0.10	< 0.10		< 0.10	0.46	< 0.10	< 0.10	< 0.10	< 0.10	< 0.10	}	< 0.10
BOD, 5-day				3.0	5.0	<1.0						6.3
Chloride	276	5 563		128	131	108	13.	195	169	163		71.5
COD	186	85.	3	< 20	65	< 20	<2	237	51.4	<20		< 20
Cyanide				<.10	< 0.10	< 0.10						< 0.10
Fluoride	0.1	7 0.1	5	1.06	1.74	1.3	0.6	3 0.82	0.65	1.32	1	0.31
Nitrate as N	1.30	6 2.7	2	0.87	0.30	1.79	0.6	0.65	0.65	0.23		< 0.10
Oil and Grease		1		< 0.10	< 0.10	< 0.10						< 0.10
pH (standard pH + lts)	7.	1 7.	5	9.3	10.4	9.	7 10.	2 9.1	8.1	10.0	)	7.6
Sulfate	109:	<del></del>		598	677	566	40.	411	660	726		412
Phenois	T	1	1		1							
Phosphorus, Total	1	1	T							L,,		
TOC	2.5	8 4.8	6	7.46	49.3	14.0	0 15.	1 10.	5.1	8.34		6.63
Total Dissolved Solids	1770	<del></del>	7	1459	2204	1855	134	2301	1590	1893		848
Turbidity (nephelometric turbidity units)	8.5	0						9,4	4 6.7	o		

Summary of Analyt Taults for Inorganic
Co Junds
at Dixon Marquette ........t Company, 1998-2000
(all values in mg/l)

979/99   11/1/99   372/00   6/23/00	MW4-WT 9/29/00  0.205 <0.001 0.002 0.122 <0.001 1.78 <0.001 0.002 0.002 0.002 0.002 0.002 0.002 0.003 0.001 51.6	MW4-WT 11/36/00 0.104 <0.001 0.004 0.190 <0.001 2.14 <0.001 155 0.005 0.001 0.009 3.80 0.002 49.7	MW4-WTF 11/30/00 0.055 <0.001 0.057 <0.001 0.232 <0.001 222 0.006 <0.001 0.007	7/26/99 6.92 <0.001 6.005 0.110 <0.001 0.09 <0.001 262 6.004 0.008 2.84	MW9-WT 979799  1.51 <0.001 0.007 0.214 <0.001 0.68 <0.001 424 0.005 0.020 0.013 7.57	MWS-VFT 11/1/99 1.65 <0.001 0.011 0.098 <0.001 0.094 <0.001 279 0.005 0.001 0.005 0.005 0.005	3/2/00 0.781 <0.001 0.006 0.031 <0.05 <0.001 203 0.004 0.007 0.025	MW9-WT 6/23/00 <0.04 <0.001 0.047 <0.001 <0.05 <0.001 104 0.002 0.002	MW9-WT 9/29/00 0.456 <0.001 0.003 0.051 <0.001 <0.05 <0.001 0.001 0.002
Aluminura         .33         0.715         1.30         <0.04	0.205 <0.001 0.002 0.122 <0.001 1.78 <0.001 164 0.002 0.002 0.008 2.23 0.001 51.6	0,104 <0.001 0,004 0,190 <0.001 2,14 <0.001 155 0,005 0,009 3,80 0,002	0.055 <0.001 <0.001 0.057 <0.001 0.232 <0.001 222 0.066 <0.001 0.007	6,92 <0.001 6,005 0,110 <0.001 0,09 <0.001 262 0,004 0,003 2,84	1.31 <0.001 0.007 0.214 <0.001 0.08 <0.001 424 0.005 0.020 0.013	1.65 <0.001 0.011 0.098 <0.001 0.094 <0.001 279 0.005 0.012 0.037	0.781 <0.001 0.006 0.031 <0.001 <0.001 205 0.004 0.007	<0.04 <0.001 0.004 0.047 <0.001 <0.05 <0.001 104 0.002	0.456 <0.001 0.003 0.051 <0.001 <0.05 <0.001 66.7 0.001 0.002
Antimony         < 0.001         < 0.001         < 0.001         < 0.001           Arsenic         0.003         0.003         0.006         0.002           Barium         0.149         0.123         0.117         0.063           Beryllium         < 0.001	<0.001 0.002 0.122 <0.001 1.78 <0.001 164 0.002 0.002 0.008 2.23 0.001 51.6	<0.001 0.004 0.190 <0.001 2.14 <0.001 155 0.005 0.001 0.009 3.80 0.002	<0.001 <0.001 <0.001 0.057 <0.001 0.232 <0.001 222 0.006 <0.001 0.007 0.953	<0.001 0.005 0.110 <0.001 0.09 <0.001 262 0.004 0.004 0.008	<0.001 0.007 0.214 <0.001 0.08 <0.001 424 0.005 0.020	<0.001 0.091 0.098 <0.001 0.094 <0.001 279 0.005 0.012 0.037	<0.001 0.006 0.081 <0.001 <0.05 <0.001 205 0.004 0.007 0.025	<0.001 0.004 0.047 <0.001 <0.05 <0.001 104 0.002 0.001	<0.001 0.003 0.051 <0.001 <0.05 <0.001 66.7 0.001
Arsenic         0.003         0.003         0.003         0.002           Barium         0.149         0.123         0.117         0.063           Beryllium         <0.001	0.002 0.122 <0.001 1.78 <0.001 164 0.002 0.002 0.008 2.23 0.001 \$1.6	0.004 0.190 <0.001 2.14 <0.001 155 0.005 0.011 0.009 3.80 0.002	<0.001 0.057 <0.001 0.232 <0.001 222 0.006 <0.001 0.007	0.005 0.110 <0.001 0.09 <0.001 262 0.004 0.004 0.008 2.84	0.007 0.214 <0.001 0.08 <0.001 424 0.005 0.020	0.011 0.098 <0.001 0.094 <0.001 279 0.005 0.012 0.037	0.006 0.081 <0.001 <0.05 <0.001 205 0.004 0.007 0.025	0.004 0.047 <0.001 <0.05 <0.001 104 0.002	0.003 0.051 <0.001 <0.05 <0.001 66.7 0.001 0.002
Barium         0.149         0.123         0.117         0.063           Beryllium         <0.001	0.122 <0.001 1.78 <0.001 164 0.002 0.002 0.008 3.23 0.001 51.6	0.190 <0.001 2.14 <0.001 155 0.005 0.011 0.009 3,80 0.002	0.057 <0.001 0.232 <0.001 222 0.066 <0.001 0.007	0.110 <0.001 0.09 <0.001 262 0.004 0.004 0.008 2.84	0.214 <0.001 0.08 <0.001 424 0.005 0.020 0.013	0.098 <0.001 0.094 <0.001 279 0.005 0.012 0.037	0.081 <0.001 <0.05 <0.001 205 0.004 0.007	0.047 <0.001 <0.05 <0.001 104 0.002 0.001	0.051 <0.001 <0.05 <0.001 66.7 0.001 0.002
Beryllium         <0.001         <0.001         <0.001         <0.001           Boron         1.05         1.21         1.53         0.11           Cadmium         <0.001	<0.001 1.78 <0.001 164 0.002 0.002 0.008 2.23 0.001 51.6	<0.001 2.14 <0.001 155 0.005 0.011 0.009 3,80 0.002	<0.001 0.232 <0.001 222 0.066 <0.001 0.007	<0.001 0.09 <0.001 262 0.004 0.004 0.008 2.84	<0.001 0.08 <0.001 424 0.005 0.020 0.013	<0.001 0.094 <0.001 279 0.005 0.012 0.037	<0.001 <0.05 <0.001 205 0.004 0.007	<0.001 <0.05 <0.001 104 0.002 0.001	<0.001 <0.05 <0.001 66.7 0.001 0.602
Boron         1.05         1.21         1.53         0.11           Cadmium         <0.001	1.78 <0.001 164 0.002 0.002 0.008 2.23 0.001 51.6	2.14 <0.001 155 0.005 0.011 0.009 3.80 0.002	0.232 <0.001 222 0.006 <0.001 0.007 0.953	0.09 <0.001 262 0.004 0.004 0.008	0.08 <0.001 424 0.005 0.020 0.013	0.094 <0.001 279 0.005 0.012 0.037	<0.05 <0.001 205 0.004 0.007 0.025	<0.05 <0.001 104 0.002 0.001	<0.05 <0.001 66.7 0.001 0.602
Cadmium         <0.001         <0.001         <0.001         <0.001         <0.001         <0.001         <0.001         <0.001         <0.001         <0.002         <0.002         <0.002         <0.002         <0.002         <0.002         <0.002         <0.002         <0.001         <0.001         <0.001         <0.001         <0.001         <0.001         <0.001         <0.001         <0.001         <0.001         <0.001         <0.001         <0.001         <0.001         <0.001         <0.001         <0.001         <0.001         <0.001         <0.001         <0.001         <0.001         <0.001         <0.001         <0.001         <0.001         <0.001         <0.002         <0.002         <0.002         <0.002         <0.002         <0.002         <0.002         <0.002         <0.002         <0.002         <0.002         <0.002         <0.003         <0.001         <0.001         <0.001         <0.001         <0.001         <0.001         <0.001         <0.001         <0.001         <0.002         <0.002         <0.002         <0.002         <0.001         <0.002         <0.002         <0.001         <0.002         <0.002         <0.001         <0.002         <0.002         <0.002         <0.003         <0.001         <0.001         <0.002	<0.001 164 0.002 0.002 0.008 2.23 0.001 51.6	<0.001 155 0.005 0.011 0.009 3,80 0.002	<0.001 222 0.006 <0.001 0.007 0.953	<0.001 262 0.004 0.004 0.008 2.84	<0.001 424 0.005 0.020 0.013	<0.001 279 0.005 0.012 0.037	<0.001 205 0.004 0.007 0.025	<0.001 104 0.002 0.001	<0.001 66.7 0.001 0.002
Calcium         175         180         261         125           Chromium         0.001         0.002         0.002         0.002           Cobalt         0.007         0.006         0.007         <0.001	164 0.002 0.002 0.008 2.23 0.001 51.6	155 0.005 0.011 0.009 3,80 0.002	222 0.066 <0.001 0.007 0.953	262 0.004 0.004 0.008 2.84	424 0.005 0.020 0.013	279 0.005 0.012 0.037	205 0.004 0.007 0.025	104 0.002 0.001	66.7 0.001 0.002
Chromium         0.001         0.002         0.002         0.002           Cobalt         0.007         0.006         0.007         <0.001	0.002 0.002 0.008 2.23 0.001 51.6	0.005 0.011 0.009 3,80 0.002	0.066 <0.001 0.007 0.953	0.004 0.004 0.008 2.84	0.005 0.020 0.013	0.005 0.012 0.037	0.004 0.007 0.025	0.002 0.001	0.001 0.602
Cobalt         0.007         0.006         0.007         <0.001           Copper         0.004         0.008         0.011         <0.001	0.002 0.008 2.23 0.001 51.6	0.011 0.009 3,80 0.002	<0.001 0.007 0.953	0.004 0.008 2.84	0.020 0.013	0.012 0.037	0.007 0.025	0.001	0.002
Copper         0.004         0.008         0.011         < 0.001           Iron         3.92         2.60         7.97         1.08           Lead         0.002         0.003         < 0.001	0.008 2.23 0.001 51.6	0.009 3,80 0.002	0.007 <i>0.95</i> 3	0.008 2.84	0.013	0.037	0.025		**********
Iron	2.23 0.001 51.6	3,80 0.002	0.953	2.84				< v.w) I	ניליט.ט
Lead         0.002         0.003         <0.001         <0.001           Magnesium         22.0         62.0         69.7         40.1           Manganese         0.368         1.67         1.07         0.171           Mercury         <0.0002	0.001 51.6	0.002			7.27		2.62	0.032	0.772
Magnesium         22.0         62.0         69.7         40.1           Manganese         0.368         1.67         1.07         0.171           Mercury         <0.0002	51.6		0.004	Λ ΩΛ <b>ξ</b> 1	0.014	0.011	< 0.001		
Manganese         0.368         1.67         1.07         0.171           Mercury         <0.0002		77.7	91.1	0.005 <b>93.</b> 4	113	91.2	75.3	< 0.001	0.004
Mercury         < 0.0002         < 0.0002         < 0.0002         < 0.0002         < 0.0000         < 0.0000         < 0.0000         < 0.0000         < 0.0003         < 0.0003         < 0.0003         < 0.0003         < 0.001         < 0.001         < 0.001         < 0.001         < 0.001         < 0.001         < 0.001         < 0.001         < 0.001         < 0.001         < 0.001         < 0.002         < 0.001         < 0.001         < 0.001         < 0.001         < 0.002         < 0.001         < 0.001         < 0.001         < 0.002         < 0.002         < 0.002         < 0.002         < 0.002         < 0.002         < 0.002         < 0.002         < 0.002         < 0.002         < 0.002         < 0.002         < 0.002         < 0.002         < 0.002         < 0.002         < 0.002         < 0.002         < 0.002         < 0.002         < 0.002         < 0.002         < 0.002         < 0.002         < 0.002         < 0.002         < 0.002         < 0.002         < 0.002         < 0.002         < 0.002         < 0.002         < 0.002         < 0.002         < 0.002         < 0.002         < 0.002         < 0.002         < 0.002         < 0.002         < 0.002         < 0.002         < 0.002         < 0.002         < 0.002         < 0.002         < 0.002         < 0.002 <td>0.510</td> <td>2.88</td> <td>0.167</td> <td>1.36</td> <td>1.66</td> <td>2.59</td> <td></td> <td>51.9</td> <td>26.5</td>	0.510	2.88	0.167	1.36	1.66	2.59		51.9	26.5
Nickel         0.025         0.015         0.012         0.003           Potassium         243         267         201         172           Selenium         0.002         0.002         <0.001	< 0.0002	< 0.0002	< 0.0002	< 0.0002	<0.0002		1.32	< 0.001	0,448
Potassium         243         267         201         172           Selenium         0.002         0.002         <0.001	0.009	0.025	0.005	0.012		< 0.0002	< 0.0002	< 0.0002	0.0004
Selenium         0.002         0.002         < 0.001         < 0.001           Silver         < 0.001	228	220	280	558	0.094 520	0.057	0.052	0.011	0.008
Silver         < 0.001         < 0.001           Sodium         55.2         25.4         23.2         13.8           Thallium         < 0.001	0.002	0.006				1900	621	527	313
Sodium         55.2         25.4         23.2         13.8           Thallium         < 0.001	0.052	0.006	0.002	0.902	0.003	0.009	0.007	< 0.001	0.003
Thallium         < 0.001         < 0.001         < 0.001         0.002           Vanadium         < 0.001	25.4	31.0		< 0.001	<0 001	< 0.001			
Vanadium <0.001 0.006 0.006 0.004		21.0	22.4	38.8	96.6	74.2	59.5	47.9	35.3
	<0.001	0.001	< 0.001	< 0.001	< 0.001	< 0.001	0.001	< 0.001	<0.00:
Zinc 0.038 0.021 0.022 <0.001	0.023	0.003	0.008	0.007	0.010	0.011	< 0.001	0.006	0.004
A!kalinity, phenol. <1.0 <1.0	<1.0	<1.0	0.008	0.032	0.068	0.050	0.028	< 0.001	0.014
Alkalinity, total 364 380 306 295.0	356.0	370		418			<1.0	<1.0	<1.0
Alkalinity, bicarbonate 345 360 297 279.0	350.0	360		391	420	540	480	500.0	550.0
Annmonia nitrogen 0.34 < 0.10 < 0.10 < 0.10	0.32	<0.10		< 0.10	400 0.28	505	398	480.0	376.0
BOD, 5-day 3.4 <1.0	- 0.32	V0.10				< 0.10	<0.10	< 0.10	0.31
Chloride 87.1 96.9 77.3 42.9	76.2	88.2		2.9 164	306	3.0 373			
COD 25 <20 <20 200	22.9	67.6		<20			219	145	191
Cyanide <0.10 <0.10		07.0			< 20	<20	<20	447	<20
Fluoride 0.23 0.24 0.16 0.28	0.29	0.31		<0.10 0.52	<0.10 0.21	< 0.10			
Nitrate as N <0.10 <0.10 <0.10 0.40	0.10	< 0.10		< 0.10	0.21		0.30	0.71	0.75
Oil and Grease <0.10 <0.10	~~~~			<0.10		1.33	0.11	< 0.10	<0.10
pH (standard pH units) 8.2 7.7 7.1 7.3	7.2	7.7			<0.10	<0.10			
Sulfate 424 487 472 365	410	432		7.8 690	8.2 986	7.8	7.5	7.9	7.6
Phenois		750		0,0	700	1000	959	506	• 717
Phosphorus, Total									
TOC 27.6 3.58 9.05 9.27	2.35	3.07		16.9	5.82	16.8	<del>;; . </del>	<del></del>	
Total Dissolved Solids 969 978 853 683	832	876		1379	1876	3120	13.1 3500	11.4	3.08
Turbidity (nephelometric turbidity units) 28.6	65.00			-23/3	10/0	3120	1300	3.40	66,20

Danamatan	MW9-WT	MW9-WTF	LW-1	LW-1	LW-1	LW-1	LW-i	LW-1	LW-1	LW-1	LW-1	LW-1	LW-1
Parameter	11/30/00	11/30/00	6/17/98	7/24/98	10/26/98	2/16/98	7/26/99	9/9/99	11/1/99	3/2/00	6/23/00	9/29/00	11/30/00
Aluminum	0.180	0.008		7.3	3.3		2.45	2.02	4.57	0.670	< 0.04	0.305	0.009
Antimony	< 0.001	< 0.001		< 0.50	< 0.20		< 0.001	< 0.001	< 0.001	< 0.001	< 0.001	< 0.001	0.001
Atsenic	0.098	0.006		< 0.400	< 0.160		0.008	0.009	0.021	0.010	0.007	0.013	0.021
Barlum	0.070	0.070		0.190	0.156		0.096	0.109	0.076	0.081	0.054	0.070	0.265
Beryllium	< 0.001	< 0.001	< 0.150	< 0.050	< 0.020		< 0.001	< 0.001	< 0.001	< 0.001	< 0.001	< 0.601	< 0.001
Boron	0.114	0.110		0.22	< 0.50		0.20	0.27	0.276	0.233	< 0.05	0.176	0.825
Cadmium	< 0.001	< 0.001		< 0.100	< 0.040		< 0.001	0.001	0.003	0.003	< 0.001	< 0.001	0.007
Calcium	117	104		620	600		758	635	673	581	397	529	623
Chromium	0.002	0.002		< 0.020	< 0.040		9.009	0.007	0.022	0.010	0.005	0.008	0.028
Cobalt	0.001	0.001	< 0.300	< 0.020	< 0.040		0.023	0.035	0.043	0.028	0.007	0.016	0.048
Copper	0.009	0.006	< 0.300	< 0.020	< 0.040		0.038	0.048	0.111	0.043	0.018	0.048	0.103
Iron	0.375	0.298	48	6.9	6.5		8.48	7.35	10.1	6.41	1.00	4.44	3.42
Lead	0.001	<del> </del>		< 0.50	< 0.020		0.015	0.016	9.019	0,002	< 0.001	0.002	< 0.001
Magnesium	59.5	60.0		220	220		287	228	30.5	216	115	168	136
Manganese	0.657	0.602	7.9	8.0	8.2		8.17	2.32	12.6	#21	3.59	5.76	29.2
Mercury	< 0.0002		0.0002	< 0.0002	< 0.00020		< 0.0002	0.0004	< 0.0002	< 0.0002	< 0.0002	0.0006	0.0034
Nickel	0.010	<del> </del>		< 0.250	< 0.100		0.043	0.055	9.082	0.044	0.017	C.037	0.079
Potassium	909	<del> </del>	5,109	4,300	5,500	5,900	6,512	6080	6074	7080	6220	6920	7420
Selenium	0.004	<del> </del>	<del> </del>	< 0.75	< 0.30		0.021	0.027	0.074	0.037	0.030	0.050	0.104
Silver	1		<u> </u>	< 0.100	< 0.040		< 0.001	< 0.001	< 0.001				
Sodium	87.1	86.4	320	280	330	320	195	366	379	396	337	387	388
Thallium	0.001	< 0.001	1	< 5.0	<2.0		< 0.001	< 0.001	0.002	6.002	0.001	0.003	0.017
Vanadium	0.005		<del></del>	< 0.250	< 0.100		0.016	0.008	0.024	0.013	0.005	0.007	0.033
Zina	0.009	0.009		< 0.100	< 0.040		0.021	0.022	0.077	0.030	< 0.001	0.028	0.029
Alkalinity, phenol.	<1.0			<5.0	<5.0				<u> </u>	<1.0	<1.0	<1.0	<1.0
Alkalinity, total	631	3	1	310	330		295	242	308	254	310.0	276.0	280
Alkalinity, bicarbonate	595		1	310	330		287	240	270	198	300.0	354.0	259
Ammonia nitrogen	< 0.10			5.2	7.7		5.61	4.53	2.31	3.66	6.09	13.7	1.40
BOD, 5-day	1	1		<5.0	< 5.0	1	2.7	4.6	24				
Chloride	293		1,600	1,900	2,000	1,700	1,445	2121	3090	2200	3096	3890	3588
СОР	23.	5		66	60	5	<20	42	< 20	25	542	257	26.5
Cyanide				< 0.0050	< 0.0050	)	< 0.10	<0.10	< 0.10				
Fluoride	1.12	2	1	0.40	0.48	3	0.31	0.44	0.22	0.82	0.86	< 0.05	0.30
Nitrate as N	0.1	7	21	1.3	12	2	< 0.10	< 0.10	< 0.10	< 0.10	< 0.10	< 0.10	< 0.10
Oil and Grease	1		1	< 1.5	3.5		2.2	< 0.10	< 0.10				
pH (standard pH units)	8.	3	7.4	7.2	7.4	7.4	7.7	8.3	7.8	7.1	7.5	7.3	7.8
Sulfate	91:		6,200	5,800	5,900	6,300	4,444	6731	6740	6400	6441	7339	7830
Phenols	1		T	< 0.029	< 0.020						<u> </u>		L
Phosphorus, Tetal	1	T	1.3				I					1	L
TOC	5.7	6		10	5 22		23.8	24.1	30.8	81.3	22.1	13.8	17.4
Total Dissolved Solids	189	4	13,000	14,000	11,000	16,000	13,200	13120	1326	14840	7115	15100	15180
Turbidity (nephelometric turbidity units)			<1								2.8	5 21.0	<u></u>

Parameter	LW-1F	LW-3	LW-3	LW-3	LW-3	LW-3	LW-3	LW-3F	SPRING	SPRING	SPRING	SPRING	SPRING	SPRING
	11/30/00	7/26/99	9/9/99	11/1/99	6/23/00	9/29/00	11/30/00	11/30/00	7/26/99	9/9/99	11/1/99	3/2/00	6/23/00	9/29/00
Aluminum	< 0.64	0.13	0.59	2.28	< 0.04	0,231	0.404	< 0.04	0.08	0.10	0.131	0.027	< 0.04	0.031
Antimony	< 0.001	< 0.001	< 0.001	< 0.001	<0.001	< 0.001	< 9.001	< 0.001	< 0,001	< 0.001	< 0.001	< 0.001	< 0.001	< 0.001
Arsenic	0.015	0.001	0.002	0.057	< 0.001	0.008	< 0.001	< 0.00;	< 0.001	< 0.001	0.002	0.001	0.007	< 0.001
Barium	0.122	0.196	0.297	0.061	0.067	0.109	0.091	0.090	0.066	0.033	0.066	0.085	0.053	0.079
Beryllium	< 0.001	< 0.001	< 0.001	< 0.001	< 0.001	< 0.001	< 0.001	< 0.001	< 0.001	< 0.001	< 0.001	< 0.001	< 0.001	< 0.001
Boron	0.375	0.09	0.13	0.284	< 0.05	0.132	• 0.159	0.169	< 0.05	< 0.05	< 0.05	< 0.05	< 0.05	< 0.05
Cadmium	0.002	<0.001	< 0.001	0.002	< 0.001	< 0.001	< 0.001	< 0.001	< 0.001	< 0.001	< 0.001	< 0.001	< 0.001	< 0.001
Calcium	485	1025	768	715.	271	309	387	286	150	73.3	143	139	112	134
Chromium	0.012	0.003	0.001	0.069	0.003	0.008	0.001	0.001	0.002	< 0.001	0.002	0.003	0.005	0.004
Cobalt	0.029	0.006	0.025	0.018	Q.002	0.007	0.003	0.004	< 0.001	< 0.001	< 0.001	< 0.001	0.007	< 0.001
Copper	0.056	0.002	0.005	0.053	<0.001	0.003	0.007	0.002	0.002	0.015	0.003	0.005	0.017	0,002
Iron	4.92	2.53	4.39	7.97	0.848	15.7	0.876	0.434	0.254	0.144	0.373	0.230	1.00	< 0.003
Lead	< 0.001	< 0.001	0.004	0.009	<0.001	0.002	0.004	0.002	< 0.001	0.006	< 0.001	<0.001	< 0.001	0.301
Magnesium	133	109	128	106	83.4	93.7	90.6	85.1	67.6	32,7	63.4	53.3	49.5	59.6
Manganese	13.0	2.61	1.93	9.19	0.830	186	1.06	1.59	0.041	0.027	9.053	0.025	3.59	0.042
Mercury	0.0013	< 0.0002	< 0.0002	< 0.0002	< 0.0002	< 0.0002	< 0.0002	< 0.0002	< 0.0002	< 0.0002	< 0.0002	< 0.0002	< 6.0002	<0.0002
Nickel	0.046	0.023	0.065	0.048	0.008	0.019	0.014	0.012	0.001	0.004	0.003	0.002	0.016	0.002
Potassium	7250	60.6	58.4	44.9	60.5	58.0	71.9	68.3	134	71.5	164	77.3	55.8	138
Selenium	0.062	0.005	0.005	0.023	< 0.001	< 0.001	0.002	< 0.001	0.002	< 0.001	0.003	0.006	0.030	0.001
Silver		<0.001	< 0.001	< 0.001					< 0.001	< 0.001	< 0.001			
Sodium	383	9.30	8.47	9.61	7.24	9.10	8.40	8.50	23.4	15.7	26.6	30.4	18.6	24.8
Thallium	0.005	0.002	0.003	< 0.001	<0.001	< 0.001	< 0.001	< 0.001	< 0.001	< 0.001	< 0.001	< 0.001	0,001	< 0.001
Vanadium	0.009	0.013	< 0.001	0.029	0.008	0.009	0.002	< 0.001	0.003	< 0.001	0.002	0.004	0.005	0.004
Zinc	0.020	0.013	0.016	0.200	< 0.001	0.918	0.018	0.026	0.005	0.069	0.012	0.010	< 0.001	0.008
Alkalinity, phenol.					<1.0	<1.0	< 1.0					<1.0	<1.0	<1.0
Alkalinity, total		367	328	564	379.0	516.0	546		254	342	354	320	318.0	340.0
Alkalinity, bicarbonate		360	310	515	570.0	332.0	497		245	330	310	297	310.0	326.0
Ammonia nitrogen		< 0.10	0.25	< 0.10	< 0.10	104	< 0.10		< 0.10	< 0.10	<0.10	< 0.10	< 0.10	< 0.10
BOD, 5-day		2.9	36.6	24					2.6	<1.0	< 1.0			
Chloride		21.0	18.1	35.1	37.3	74.7	62.7		82.8	95.9	114	98.4	<b>\$4.</b> 6	99.8
COD		<20	45	<20	<20	51.4	<20		<20	<20	<20	< 20	421	<20
Cyanide	<u> </u>	<0.10	< 0.10	< 0.10					< 0.10	< 0.10	< 0.10			
Fluoride		0.21	0.20	0.20	0.26	0.22	0.25		0.21	0.16	0.17	0.13	0.17	0.17
Nitrate as N		< 0.10	< 0.10	< 0.10	< 0.10	0.16	0.41	<u> </u>	11.3	13.0	11.1	4.75	10.8	9.42
Oil and Grease		< 0.10	<0.10	< 0.10			,		< 0.10	< 0.10	<0.10			
pH (standard pH units)		7.3	7.7	7.4	6.8	6.8	8.2		8.1	8.3	7.7	7.0	7.3	7.1
Sulfate		429	416	403	431	474	471		210	235	503	198	137	255
Phenois								<u> </u>			ļ			
Phosphorus, Total	<u></u>		i	-					<u> </u>					
TOC		12.8	9.36	11.8	13.4	2.57	3.94	<u></u> ,	10.7	8.48	5.06	6.00	6.70	1.17
Total Dissolved Solids		783	858	867	867	860	898	<u> </u>	682	730	825	638	533	722
Turbidity (nephelometric turbidity units)					17.2								0.15	0.15

Parzmeter	SPRING 11/30/00	RRUS 3/2/00	RRUS 6/23/00	RRUS 9/19/00	RRUS 11/30/00	RRDS 3/2/00	RRDS 6/23/00	RRDS 9/29/00	RRDS 11/30/00
Aluminum	< 0.04	0.191	0.179	0.435	< 0.04	0.172	0.163	0.535	< 0.04
Antimony	< 0.001	< 0.001	< 0.001	< 0.001	<0.001	< 0.001	< 0.001	< 0.001	< 0.001
Arsenic	< 0.001	0.002	< 0.001	0.001	< 0.001	0.002	< 0.001	< 0.001	< 0.001
Barium	0.065	0.064	0.064	0.078	0.046	0.063	0.063	0.079	0.041
Beryllium	< 0.001	< 0.001	< 0.001	< 0.001	< 0.001	< 0.001	< 0.001	< 0.001	< 0.001
Boron	0.035	< 0.05	< 0.05	< 0.05	0.042	< 0.05	< 0.05	< 0.05	0.036
Cadmium	< 0.001	< 0.001	< 0.001	< 0.001	< 0.001	< 0.001	< 0.001	< 0.001	< 0.001
Calcium	135	66.1	66.6	71.8	75.8	65.8	64.3	74.1	73.0
Chromium	< 0.003	0.002	0.001	0.001	< 0.003	0.002	0.002	0.001	< 0.003
Cobalt	< 0.001	< 0.001	< 0.001	0.002	< 0.001	< 0.001	< 0.001	< 0.001	< 0.001
Copper	0.002	0.005	0.001	0.005	< 0.001	0.005	< 0.001	0.004	0.001
Itoa	0.072	0.463	0.481	0.800	0.046	0.345	0.476	0.947	0.038
Lead	<0.001	< 0.001	0.002	0.002	<0.001	< 0.001	0.002	0.003	< 0.001
Magnesium	59.6	32.2	31.6	37.3	40.9	32.0	31.4	38,6	41.2
Manganese	0.033	0.068	0.055	0.146	0.025	0.063	0.044	0.153	0.04
Mercury	< 0.0002	< 0.0002	< 0.0002	< 0.0002	< 0.0002	< 0.0002	< 0.0002	< 0.0002	< 0.0002
Nickel	0.001	0.003	0.006	0.002	0.001	0.002	0.001	0.002	0.001
Potassium	178	5.55	7.00	6.70	3.20	5.13	7.23	3.80	2.80
Selenium	< 0.001	0.008	< 0.001	0.001	< 0.001	0.005	< 0.001	0.002	< 0.001
Silver									
Sodium	25.2	20.2	11.3	18.2	20.5	20.0	12.4	18.7	21.0
Thallium	< 0.001	< 0.001	< 0.001	< 0.001	< 0.001	< 0.001	< 0.001	< 0.001	< 0.001
Vanadium	< 0.001	0.004	0.006	0.005	< 0.001	0.004	0.006	0.005	< 0.001
Zinc	< 0.001	0.021	< 0.001	0.025	< 0.001	0.014	< 0.001	0.019	< 0.001
Alkalinity, phenol.	< 1.0	<1.0	<1.0	16.0	16	<1.0	<1.0	<1.0	16
Alkalinity, total	338	200	210.0	242.0	272	197	213.0	230,0	270
Alkalinity, bicarbonate	319	139	185.0	187.0	231	109	207.0	216.0	227
Ammonia nitrogen	< 0.10	<0.1J	< 0.10	0.37	<0.10	< 0.10	< 0.10	0.13	< 0.10
BOD, 5-day									
Chloride	112	41.4	28.6	39.5	44.5	40.6	28.9	46.3	50.1
COD	< 20	< 20	<20	31.4	29.4	<20	126	<20	<20
Cyanide									
Fluoride	0.21	0.14	0.14	0.14	0.15	0.18	0.14	0.13	0.14
Nitrate as N	8.75	5.69	5.76	4.27	4.55	5.71	5.85	4.12	3.85
Oii and Grease				L					
pH (standard pH units)	7.7	7.6	8.1	8.1	8.7	7.8	8.1	8.0	8.7
Sulfate	320	30.3	29.1	33.6	39.9	30.4	28.0	35.3	33.5
Phenois	ļ								
Phosphorus, Total						ļ			<u> </u>
TOC	1.95	9.94	9.14	3.87	4.67	9.25	10.9	4.00	4.47
Total Dissolved Solids	805	304	280	308	360	306	281	314	357
Turbidity (nephelometric turbidity units)		ŀ	33.00	45.00	ļ		39.60	45.00	i

Parameter	MW1-S	MW1-S	MW1-S	MW1-S	MW1-S	MW1-S	MW2-S	MW2-S	MW2-S	MW3-S	MW3-S	MW3-S
	7/24/98	10/26/98	2/16/99	7/26/99	9/9/99	11/1/99	7/26/99	9/9/99	11/1/99	7/26/99	9/9/99	11/1/99
Acenaphthene	<21	< 10										
Acenaphthylene	<21	< 10										
Alachier (lasso)	<0.5	< 0.5	<0.5	< 1.0	<1.0	<1.0	< 1.0	< 1.0	< 1.0	<1.0	<1.0	<1.0
Atrazine Anthracene	<0.5	<0.5 <10	<0.5	< 1.0	<1.0	<1.0	<1.0	< 1.0	< 1.0	<1.0	< 1.0	<1.0
Benzidine	<21 <210	<100										
Benzo(a)anthracene	<21	<10										
Benzo(b) fluoranthene	<21	<10										
Benzo(k)fluoranthene	<21	<10										
Benzo(a)pyrene	<21	<10							7			
Benzo(g,h,i)perylene	<21	<10										
Benzyl alcohol	<21	< 10										
Benzy' butyl phthalate	<21	< 10			`							
Bis(2-chloroethyl) ether	<21	< 10										
Bis(2-chlorethoxy)methane	<21	<10										
Bis(2-ethylhexyi)phthalate	<21	<10		<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0
Bis(chloromethyl)ether				< 1.0	< 1.0	<1.0	<1.0	< 1.0	<1.0	<1.0	<1.0	< 1.0
Bis(2-chloroisopropyl)ether	<21					ļ	ļ	<u> </u>				
4-Bromophenyl phenyl ether	<21	·										
4-Chloroaniline	<21			ļ	ļ	<b></b>				<b> </b> -		
2-Chloronaphthalene	<21					<del> </del>						
4-Chlorophenylphenyl ether	<21	<10		<del> </del>		<del> </del>				<b> </b>	ļ	
Chrysene Dibenzo(a,h) anthracene	<21 <21	<10 <10				<b></b>	ļ	ļ		<b></b>		
Di-n-butylphthalate	<21			< 1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.6	<1.0
1.2-Dichlorobenzene	<21					\$1.0	<u> </u>		<u> </u>	\ <u>````</u>	\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\	
1,3-Dichlorobenzene	<21											
1,4-Dichlorobenzene	<21	< 10	<del></del>							<b> </b>		
3,3-Dichlorobenzidine	<100	<50		1				<del></del>				
Diethyl phthafate	<21	<10		<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0
Dimethyl phthalate	<21	<10		< 1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0
2,4-Dinitrotolucne	<21	< 10				<u> </u>						
2,6-Dinitrotoluene	<21	< 10				<u> </u>	ļ <u>.</u>		ļ		ļ	ļ
Di-n-octylphthalate	<21	< 10		ļ	ļ	<u> </u>	ļ	ļ		ļ	ļ	
Fluoranthene	<21	< 10	<b> </b>	<del> </del>		ļ		ļ	ļ	<u> </u>	ļ	ļ
Fluorene	<21	<10	<b> </b> -	<del> </del>	<del></del>	<del> </del>		<b> </b>		<del> </del> -	ļ	ļ
Hexachlero-1,3-hutadiene	<21	<10	< 10	<b></b>	<del> </del>	<b> </b>	ļ	ļ	<b> </b>	<b></b>	<del> </del>	
Hexachlorocyclopentadiene Hexachloroethane	<42 <21			ļ			ļ <del></del>		<b> </b>	<b></b>		
Indeno(1,2,3-cd)pyrene	<21		<del> </del>	<del> </del> -		<del> </del>	<del></del>				<del> </del>	<del></del>
Isophorone	<21	< 10	<del></del>	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0
2-Methylnaphthalene	<21	<del> </del>	<del></del> -	1 3.55		1	<del> </del>	<u>````</u>		<u></u>	<del>```</del>	<u>&gt;</u>
Naphthalene	<21	<del></del>	<del></del>	<1.0	< 1.0	<1.0	<1.0	<1.0	<1.0	<1.0	< 1.0	<1.0
2-Nitroaniline	<21	< 10										
3-Nitroaniline	<21	< 10										
4-Nitroaniline	<21	<10										
Nitrobenzene	<21	< 10		<u> </u>	<u> </u>							
N-Nitrosodimethylamine	<21	< 10			ļ	ļ	ļ	ļ				
N-Nitrosodiphenylamine	<21				<b> </b>	ļ		<u> </u>		ļ		<u></u>
N-Nitrosodi-n-propylamine	<21		·	<b> </b>	<b> </b>	ļ		<b> </b>	ļ	ļ	<u> </u>	<b></b>
Phenanthrene	<21	<del> </del>		<b></b>	<b> </b>		<del> </del>	ļ	<b> </b>	<b> </b>		
Pyrene	<21			<del> </del>		<del> </del>	<del> </del>	ļ	ļ	<del> </del> -	<del> </del>	ļ
Pyridine	<21			<del> </del>	<del> </del>	<del> </del>	<del> </del>	<del> </del>	<del> </del>	<del> </del>	<del> </del>	
1,2,4-Trichlorobenzene	<21	·		1	<del> </del>	<del> </del>	<del> </del>	<del> </del>	<del> </del>	<del> </del>	<del> </del>	<del> </del>
Benzoic Acid	<42			<del> </del>	<del></del>	<del> </del>	<del> </del>	<del> </del>	<del> </del>	<del> </del>	<del> </del>	<del> </del>
4-Chloro-3-methylphenol	<21	<del> </del>	<del> </del>	<del> </del>	<del> </del> -	<del> </del>	<del> </del>	<del> </del>	<del> </del>	<del> </del>	<del> </del>	ļ
2-Chlorophenol	< 21	<10	<del> </del>		<del> </del>	ļ		<del> </del>		<del> </del>	<del> </del>	<del> </del>
p-Cresol Cresole Total	<21	< 10	<del> </del>	<1.0	< 1.0	<1.0	<1.0	<1.0	<1.0	<1.0	< 1.0	<1.0
Cresols, Total  2,4-Dichlorophenol	<21			<del>                                     </del>	<del> </del>	<del> </del>	<del> </del>	<del> </del>	<del> </del>	<del> </del>	<del> </del> -	<del> </del>
2,4-Dimethylphenol	<21	<del> </del>		<del>                                     </del>	<del> </del>	<del>                                     </del>	<b> </b>	<del> </del>	<del>                                     </del>	<del> </del>	<del> </del>	<del> </del> -
2,4-Dinitrophenol	<21			<b>├</b> ──	h	1	t	t		<del> </del>	1	<del> </del>
C. C. Milaskinski				L		<del></del>	<del></del>	J.,	<del></del>	ــــــــــــــــــــــــــــــــــــــ		

Parameter	MW1-S 7/24/98	MW1-S 10/26/98	MW1-S 2/16/99	MW1-S 7/26/99	MW1-S 9/9/99	MW1-S 11/1/99	MW2-S 7/26/99	M1\V2-S 9/9/99	MW2-S 11/1/99	MW3-S 7/26/99	MW3-S 9/9/99	MW3-S 11/1/99
2-Methyl-4,6-dinitrophenol	<21	< 10				,						
2-Nitrophenol	<21	< 10										
4-Nitrophenol	<21	< 10						-				
Parathion	< 0.5	< 0.5		<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0
Pentachlorophenol	<21	< 10	<10	<1.0	<1.0	< 1.0	<1.0	<1.0	<1.0	<1.0	< 1.0	<1.0
Phenol	<21	<10		<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0
2,4,5-Trichlorophenol	<21	<10									1	<u></u>
2,4,6-Trichlorophenol	<21	< 10										
Aldrin	< 0.05	<0.10		< 0.034	< 0.034	< 0.034	< 0.034	< 0.034	< 0.034	< 0.034	< 0.034	< 0.034
alpha-BHC	< 0.05	< 0.10									30,000	
beta-BHC	< 0.05	< 0.10										
gamma-BHC	< 0.05	< 0.10	< 0.10	< 0.025	< 0.025	< 0.025	<0.025	< 0.025	< 0.025	< 0.025	< 0.025	< 0.025
delta-BHC	< 0.05	< 0.10						10.025	70,020	- 10.000	70.023	(0.02
Chlordane	<0.2	<1.0	<1.0	< 0.037	< 0.037	< 0.037	< 0.037	< 0.037	< 0.037	< 0.037	< 0.037	< 0.037
4,4'-DDD	< 0.05	<0.10	- \v	10.037		\U.U3/	(0.037	C0.037	C 0.037	CC.037	₹0.037	<0.037
4,4'-DDE	< 0.05	<0.10	J									
4,4'-DDT	< 0.05	<0.20		< 0.091	< 0.091	<0.091	< 0.091	< 0.091	< 0.091	< 0.091	<0.001	-C 001
Dieldrin	< 0.05	<0.10		<0.044	< 0.044	< 0.044	< 0.044	< 0.044	<0.091 <0.044	< 0.044	< 0.091	< 0.091
Endosulfan I	<0.05	< 0.05		<u> </u>	<u> </u>	C0.044	<b>\(\tau_{0.044}\)</b>	< 0.044	< U.U44	₹0.044	< 0.044	< 0.044
Endosulfan II	<0.05	< 0.20										
Endosulfan Sulfate	< 0.05	<0.10										
Endrin	< 0.05	< 0.10	< 9.10	< 0.039	< 0.039	< 0.020	<0.039	10.020	- O 036			
Endrin Aldehyde	<0.05	< 0.10	< 0.10	(0.039	< 0.039	< 0.039	< 0.039	< 0.039	< 0.039	< 0.039	< 0.039	< 0.039
Heptachlor	< 0.05	<0.10	<0.10	< 0.040				.0.040				·
Heptachlor Epoxide	< 0.05	< 0.10	<0.10	< 0.032	< 0.040	< 0.040	< 0.040	< 0.040	< 0.040	< 0.040	< 0.040	< 0.040
Methoxychlor	< 0.05				< 0.032	< 0.032	< 0.032	< 0.032	< 0.032	< 0.032	< 0.032	< 0.032
Toxaphene		<1.0	<1.0 <2.0	< 0.176	< 0.176	< 0.176	< 0.176	< 0.176	< 0.176	< 0.176	< 0.176	< 0.176
Aroclor-1016	< i,0	< 2.0	< 2.0	< 0.086	< 0.086	< 0.086	< 0.086	< 0.086	< 0.086	< 0.086	< 0.086	< 0.086
Aroclor-1221				< 0.050	< 0.050	< 0.050	< 0.050	< 0.050	< 0.050	< 0.050	< 0.050	< 0.050
Aroclor-1232				< 0.054	< 0.054	< 0.054	< 0.054	< 0.054	< 0.054	< 0.054	< 0.054	< 0.054
Aroclor-1242				< 0.065	< 0.065	< 0.065	< 0.065	< 0.065	< 0.065	< 0.065	< 0.065	< 0.065
Aroclor-1248				< 0.065	< 0.065	< 0.065	< 0.065	< 0.065	< 0.065	< 0.065	< 0.065	< 0.065
Aroclor-1254	<del></del>			< 0.090	< 0.090	< 0.090	< 0.090	< 0.090	<0.090	< 0.090	< 0.090	< 0.090
Aroclor-1260				< 0.100	< 0.100	< 0.100	< 0.100	< 0.100	< 0.100	< 0.100	< 0.100	< 0.100
Aroclor-1268				< 0.100	< 0.100	< 0.100	< 0.100	< 0.100	< 0.100	< 0.100	< 0.100	< 0.100
PCB-1016/1242	< 0.05	<1.0		< 0.100	< 0.100	< 0.100	< 0.100	< 0.100	< 0.100	< 0.100	< 0.100	< 0.100
PCB-1010/1242	< 0.05		<1.0									
PCB-1232	< 0.05	<1.0 <1.0	<1.0									
PCB-1248	< 0.05	<1.0 <1.0	<1.0									
PCB-1254	< 0.05		<1.0				i		———-{			
PCB-1260	< 0.05	<1.0 <1.0	<1.0 <1.0		·							
PCB-1268	~ < 0.03	< 1.0		<del></del>								
2,4-D	<2.0	< 2.0	<1.0 <2.0	< 0.29								
2,4-0 Silvex(2,4,5-TP)	<2.0			<0.29	< 0.29	< 0.29	< 0.29	< 0.29	< 0.29	< 0.29	< 0.29	< 0.29
· · · · · · · · · · · · · · · · · · ·		< 2.0	<2.0		<0.34	< 0.34	< 0.34	< 0.34	< 0.34	< 0.34	< 0.34	< 0.34
Aldicarb	< 10	<10	< 10	< 0.50	<.50	< .50	< 0.50	<.50	<.50	< 0.50	< .50	<.50
Carbofuran	< 10	< 10	< 10	< 5.0	< 5.0	< 5.0	< 5.0	< 5.0	< 5.0	<5.0	< 5.0	< 5.0
1,2-Dibromo-3-Chloropropare				< 0.10	< 0.10	< 0.10	< 0.10	< 0.10	< 0.10	< 0.10	< 0.10	< 0.10
Ethylene dibromide				< 0.10	< 0.10	< 0.10	< 0.10	< 0.10	< 0.10	< 0.10	< 0.10	< 0.10
Trifluralin			< 5									ļ

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Parameter	MW4-S	MW4-S	MW4-S	MW5-S	MW5-S	MW5-S	MW6-S	MW6-S	MW6-S	MW7-S	MW7-S	MW7-S
Acenaphthene	7/26/99	9/9/99	11/1/99	7/26/99	9/9/99	11/1/99	7/26/99	9/9/99	11/1/99	7/26/99	9/9/99	11/1/99
Acenaphthylene												
Alachlor (lasso)	<1.0	<1.0	< 1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0
Atrazine	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0
Anthracene		******	<u> </u>	71.0	<u></u>	<u></u>	~ 1.0		<u> </u>	7.1.0		
Benzidine	<b> </b>											
Benzo(a)anthracene												
Benzc(b)fluoranthene												
Benzo(k)fluoranthene												
Benzo(a)pyrene				-								
Benzo(g,h,i)perylene												
Benzyl alcohol												
Benzyl butyl phthalate	1											
Bis(2-chloroethyl) ether												
Bis(2-chlorethoxy)methane												
Bis(2-ethylhexy!)phthalate	< 1.0	<1.0	<1.0	< 1.0	<1.0	<1.0	<1.0	< 1.0	<1.0	<1.0	<1.0	<1.0
Bis(chloromethyl)ether	<1.0	<1.0		<1.0	<1.0	< 1.0	< 1.0	<1.0	<1.0	<1.0	<1.0	<1.0
Bis(2-chloroisopropyl)ethe:											````	
4-Bromophenyl phenyl ether	1		·		<del></del>							
4-Chloroaniline												
2-Chloronaphthalene						***********				<del></del> -		
4-Chlorophenylphenyl ether												
Chrysene						¥						
Dibenzo(a,h) anthracene		-										
Di-n-butylphthalate	<1.0	<1.0	<1.0	<1.0	<1.0	< 1.0	<1.0	< 1.0	< 1.0	<1.0	<1.0	<1.0
1,2-Dichlorobenzene	1			•		>1.0	~ ~ ~ ~ ~ ~ ~ ~ ~ ~ ~ ~ ~ ~ ~ ~ ~ ~ ~ ~	<u></u>	<u> </u>	- 51.0	<u></u>	
1,3-Dichlorobenzene												
1,4-Dichlorobenzene												
3,3-Dichlorobenzidine												
Diethyl phthalate	<1.0	<1.0	< 1.0	<1.0	<1.0	<1.0	< 1.0	<1.0	< 1.0	<1.0	<1.0	<1.0
Dimethyl phthalate	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0
2,4-Dinitrotoluene			1		77.0	47.10						
2,6-Dinitrotoluene												
Di-n-octylphthalate	<del></del>							<del>-</del>				
Fluoranthene											<del></del>	
Fluorene												
Hexachloro-1,3-butadiene												
Hexachlorocyclopentadiene												
Hexachloroethane												
Indeno(1,2,3-cd)pyrene												
Isophorone	< 1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	< 1.0	<1.0	<1.0
2-Methylnaphthalene											- 1,15	
Naphthalene	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	< 1.0	< 1.0	< 1.0	<1.0	<1.0	<1.0
2-Nitroaniline												
3-Nitroaniline												
4-Nitroaniline												
Nitrobenzene												
N-Nitrosodimethylamine				_								
N-Nitrosodiphenylamine												
N-Nitrosodi-n-propylamine												
Phenanthrene												
Pyrene									·			
Pyridine												
1,2,4-Trichlorobenzene												
Benzoic Acid												
4-Chloro-3-methylphenol												
2-Chlorophenol												
p-Cresol	<1.0	< 1.0	< 1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	< 1.0		
Cresols, Total	<del></del>					73,0		<u> </u>	\ <u>&gt;1.0</u>	71.0	<1.0	<1.0
2,4-Dichlorophenol	<u>-</u>						<del></del>					
2,4-Dimethylphenol										<b></b>		
2,4-Dinitrophenol	··········						<del></del>					
	I				لــــــل		L		L	L	<u> </u>	

Parameter	MW4-S 7/26/99	MW4-S 9/9/99	MW4-S 11/1/99	MW5-S 7/26/99	MW5-S 9/9/99	MW5-S 11/1/99	MW6-S 7/26/99	MW6-S 9/9/99	MW6-S 11/1/99	MW7-S 7/26/99	MW7-S 9/9/99	MW7-S 11/1/99
2-Methyl-4,6-dinitrophenol												
2-Nitrophenol												
4-Nitrophenol												
Parathion	< 1.0	<1.0	< 1.0	< 1.0	<1.0	<1.0	< 1.0	<1.0	<1.0	<1.0	<u>&lt;1.0'</u>	< 1.0
Pentachlorophenol	< 1.0	<1.0	<1.0	< 1.0	<1.0	< 1.0	<1.0	<1.0	<1.0	<1.0	< 1.0	<1.0
Pheno!	<1.0	<1.0	< 1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	< 1.0
2,4,5-Trichlorophenol												
2,4,6-Trichlorophenol												
Aldrin	< 0.034	< 0.034	< 0.034	< 0.034	< 0.034	< 0.034	< 0.034	< 0.034	< 0.034	< 0.034	< 0.034	< 0.034
alpha-BHC												
beta-BHC												
gamma-BHC	< 0.025	< 0.025	< 0.025	< 0.025	< 0.025	< 0.025	< 0.025	< 0.025	< 0.025	< 0.025	< 0.025	< 0.025
delta-BHC												
Chlordane	< 0.037	< 0.037	< 0.037	< 0.037	< 0.037	< 0.037	< 0.037	< 0.037	< 0.037	< 0.037	< 0.037	< 0.037
4,4'-DDD												
4,4'-DDE												
4,4'-DDT	< 0.091	< 0.091	< 0.091	< 0.091	< 0.091	< 0.091	< 0.091	< 0.091	< 0.091	< 0.091	< 0.091	< 0.091
Dieldrin	< 0.044	< 0.044	< 0.044	< 0.014	< 0.044	< 0.044	< 0.014	< 0.044	< 0.044	< 0.044	< 0.044	< 0.044
Endosulfan I												
Endosulfan II								<u> </u>				
Endosulfan Sulfate								<u></u>				
Endrin	< 0.039	< 0.039	< 0.039	< 0.039	< 0.039	< 0.039	< 0.039	< 0.039	< 0.039	< 0.039	< 0.039	< 0.039
Endrin Aldehyde								<u> </u>	ļ			
Heptachlor	< 0.040	< 0.040	< 0.040	< 0.040	< 0.040	< 0.040	< 0.010	< 0.040	< 0.040		< 0.040	< 0.040
Heptachlor Epoylde	< 0.032	< 0.032	< 0.032	< 0.032	< 0.032	< 0.032	< 0.032	< 0.032	< 0.032	< 0.032	< 0.032	< 0.032
Methoxychlor	< 0.176	< 0.176	< 0.176	< 0.176	< 0.176	< 0.176	< 0.176	< 0.176	< 0.176	< 0.176	< 0.176	< 0.176
Toxaphene	< 0.086	< 0.086	< 0.086	< 0.086	< 0.086	< 0.086	< 0.086	< 0.086	< 0.086	< 0.086	< 0.086	< 0.086
Arocior-1016	< 0.050	< 0.050	<0.050	< 0.050	< 5.050	< 0.050	< 0.050	< 0.050	< 0.050	+	< 0.050	< 0.050
Aroclor-1221	< 0.054	< 0.05	< 0.054	< 0.054	< 0.054	< 0.054	< 0.054	< 0.054	< 0.054	<del> </del>	< 0.054	< 0.054
Aroclor-1232	< 0.065	< 0.065	< 0.065	< 0.065	< 0.065	< 0.065	< 0.065	< 0.065	< 0.065		< 0.065	< 0.065
Aroclor-1242	< 0.065	< 0.065	< 0.065	< 0.065	< 0.065	< 0.065	< 0.065	< 0.065	< 0.065	< 0.065	< 0.065	< 0.065
Aroclor-1248	< 0.090	< 0.090	< 0.090	< 0.090	< 0.090	< 0.090	< 0.090	< 0.090	< 0.090	< 0.090	< 0.090	< 0.090
Aroclor-1254	< 0.100	< 0.100	< 0.100	< 0.100	< 0.100	< 0.100	< 0.100	< 0.100	< 0.100	< 0.100	< 0.100	< 0.100
Aroclor-1260	< 0.100	< 0.100	< 0.100	< 0.100	< 0.100	< 0.100	< 0.100	< 0.100	< 0.100	< 0.100	< 0.100	< 0.100
Aroclor-1268	< 0.100	< 0.100	< 0.100	< 0.100	< 0.100	< 0.100	< 0.100	< 0.10	< 0.100	< 0.100	< 0.100	< 0.100
PCB-1016/1242				<u> </u>		<u> </u>	<u> </u>	ļ		.	ļ	ļ
PCB-1221					<u> </u>		<u> </u>	<u> </u>	<u> </u>			
PCB-1232		1	<u> </u>	<u> </u>		<u> </u>	<b>_</b>	<u> </u>		<b></b>		ļ
PCB-1248		<u> </u>	<u> </u>	<u> </u>		<b></b>	<u> </u>	ļ	ļ	<b>_</b>	ļ	
PCB-1254		<u> </u>		<u> </u>	<u> </u>	<u> </u>		<del> </del>	ļ	ļ	·	
PCB-1260		<u> </u>	<u> </u>		<u> </u>	ļ	ļ	ļ	<u> </u>	<b></b>	<del> </del> -	ļ
PCB-1268 •			<b></b>		1	<del> </del>	<b></b>	ļ	<b>↓</b>	<del> </del>	<u> </u>	ļi
2,4·D	< 0.29	< 0.2	9 < 0.29	< 0.2	9 < 0.29	<0.2		+				1
Silvex(2,4,5-TP)	< 0.3	< 0.3	4 < 0.3	< 0.3		<del></del>					<del></del>	1
Aldicarb	< 0.50	<.5	0 <.5	< 0.50							<del></del>	
Carbofuran	< 5.0	< 5.	0 <5.0			0 < 5.					<del></del>	
1,2-Dibromo-3-Chloropropane	<.0.10	<0.1	0 <0.10	0.1	0 <0.1	1.0>			<del></del>			
Ethylene dibromide	< 0.10	0 < 0.1	0 < 0.1	<0.1	0 < 0.1	0 < 0.1	0 < 0.1	0 < 0.1	0 <0.1	0 <0.3	0 < 0.1	< 0.10
Trifluralin				<u></u>	1						1	1

Parameter	MW8-S 7/26/99	MW8-S 9/9/99	MW8-S 11/1/99	MW9-S 7/26/99	MW9-S 9/9/99	MW9-S 11/1/99	MW1-D 7/24/98	MW1-D 10/26/98	MWI-D 2/16/99	MW1-D 7/26/99	M2W1-D 9/9/99	MWI-D 11/1/99
Acenaphthene							<19	<10				
Acenaphthylene							< 19	<10		<del> </del> -		<del></del>
Alachior (lasso)	<1.0	<1.0	<1.0	<1.0	<1.0	< 1.0	<0.5	<0.5		<1.0	<1.0	<1.0
Atrazine	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.5	<0.5		<1.0	<1.0	<1.0
Anthracene					2313		<19	<10			1.11	
Benzidine							< 190	<100				
Benzo(a)anthracene							< 19	< 10				
Benzo(b)fluoranthene							<19	<10				<del>~~~~</del>
Benzo(k)fluoranthene							< 19	< 10				
Benzo(a)pyrene							< 19	< 10				
Benzo(g,h,i)perylene							< 19	< 10				
Benzyl alcohol							< 19	< 10				
Benzyl butyl phthalate							< 19	< 10				
Bis(2-chloroethyl) ether							< 19	< 10				
Bis(2-chlorethoxy)methane							< 19	< 10				
Bis(2-ethylhexyl)phthalate	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	< 19	<10		<1.0	<1.0	<1.0
Bis(chloromethyl)ether	<1.0	<1.0	<1.0	<1.0	< 1.0	<1.0				<1.0	<1.0	<1.0
Bis(2-chloroisopropyl)ether							< 19	< 10				
4-Bromophenyl phenyl ether							< 19	< 10				
4-Chloroaniline							<19	<10				. 25 - 1 - 1 - 1 - 1 - 1 - 1 - 1 - 1 - 1 -
2-Chloronaphthalene							< 19	<10				
4-Chlorophenylphenyl ether							< 19	<10				
Chrysene					~		<19	<10				
Dibenzo(a,h) anthracene							< 19	<10				
Di-n-outylphthalate	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	< 19	<10		< 1.0	<1.0	<1.0
1,2-Dichlorobenzene						1,,,	< 19	<10	<del></del>			
1,3-Dichlerobenzene							< 19	< 10				<del>~~~~~~</del>
1,4-Dichlorobenzene							<19	< 10			<del></del>	
3,3-Dichlorobenzidine							< 95	< 50				-
Diethyl phthalate	<1.0	< 1.0	<1.0	<1.0	<1.0	<1.0	<19	< 10		< 1.0	<1.0	<1.0
Dimethyl phthalate	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	< 19	< 10		< 1.0	<1.0	<1.0
2,4-Dinitrotoluene							< 19	< 10		<u></u>	71.0	
2,6-Dinitrotoluene							< 19	< 10				
Di-n-octylphthalate			•				< 19	< 10				<del></del>
Fluoranthene							<19	<10				
Fluorene							< 19	< 10				
Hexachloro-1,3-butadiene							< 19	< 10	<1(			
Hexachlorocyclopentadiene							< 38	< 20				
Hexachloroethane							< 19	< 10				
Indeno(1,2,3-cd)pyrene							< 19	< 10				
Isophorone	< 1.0	< 1.0	<1.0	< 1.0	< 1.0	< 1.0	< 19	< 10		<1.0	<1.0	<1.0
2-Methylnaphthalene							< 19	< 10			3,00	
Naphthalese	<1.0	<1.0	< 1.0	< 1.0	<1.0	<1.0	< 19	< 10	< 10	<1.0	<1.0	< 1.0
2-Nitroaniline							< 19	< 10				
3-Nitroaniline	1						< 19	< 10				
4-Nitroaniline							< 19	₹10	*			
Nitrobenzene							< 19	< 10				
N-Nitrosodimethylamine							< 19	< 10				
N-Nitrosodiphenylamine							< 19	< 10		***		
N-Nitrosodi-n-propylamine							< 19	<10				
Phenanthrene							< 19	< 10				
Pyrene							< 19	< 10				~
Pyridine							< 19	<10				
1,2,4-Trichlorobenzene					i		<19	<10				
Benzok Acid	<u> </u>				i		< 38					
4-Chioro-3-methylphenol						<del></del>		<10			<del>  </del>	
2-Chlorophenol	1						<u>&gt; 17</u> < 19	<10			<del></del> }	
p-Cresul	<1.0	<1.0	<1.0	<1.0	<del> </del>			<u> </u>		<u> </u>		
Cresols, Total	<del>  </del>	~ ~ 1.0	×1.0	~ 1.0	<1.0	<1.0				< 1.0	<1.0	< 1.0
2,4-Dichlorophenol	+						< 19	< 10			<b></b>	
2,4-Dimethylphenol							< 19	<10			<b>  </b>	
2,4-Dinitrophenol	<del>-  </del>						< 19	< 10				
s <sub>1</sub> T- 22 (IIII V) (III III)				1	L		< 19	< 10		L	<u></u>	1

Parameter	MW8-S 7/26/99	MW8-S 9/9/99	MW8-S 11/1/99	MW9-S 7/26/99	MW9-S 9/9/99	MW9-S 11/1/99	MW1-D 7/24/98	MW1-D 10/26/98	MW1-D 2/16/99	MW1-D 7/26/99	MW1-D 9/9/59	MW1-D 11/1/99
2-Methyl-4,6-dinitrophenol							<19	<10				
2-Nitrophezol							<19	<10				
4-Nitrophenol							<19	<10				
Parathion	<1.0	<1.0	<1.0	<1.0	-51.0	<1.0	<0.5	<0.5		<1.0	<1.0	<1.0
Pentachlorophenol	<1.0	<1.0	< 1.0	< 1.0	< 1.0	<1.0	< 19	< 10	<10	<1.0	<1.0	<1.0
Pheno!	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	< 19	< 10		< 1.0	<1.0	<1.0
2,4,5-Trichiorophenol							< 19	<10			<del>``</del>	
2,4,6-Trichlorophenol							<19	<10			·	
Aldrin	< 0.934	< 0.034	< 0.034	< 0.034	< 0.034	< 0.034	< 0.05	<0.10		< 0.034	< 0.034	< 0.034
alpha-BHC					10,00	10.03	< 0.05	<0.10			10.054	70.057
beta-BFC							< 0.05	< 0.10				
gainma-BHC	< 0.025	< 0.025	< 0.025	<0.025	< 0.025	< 0.025	< 0.05	<0.10	< 0.10	< 0.025	< 0.025	< 0.025
delta-BHC		40.000	10.00	10.020		50.023	< 0.05	<0.10	V0.10	₹0.025	< 0.023	C0.025
Chlordane	< 0.037	< 0.037	< 0.037	< 0.037	< 0.037	< 0.037	<0.2	<1.0	<1.0	< 0.037	< 0.037	< 0.037
4,4'-DDD	1 1000	10.057	×0.037	10.051	(0.037	V0.037	< 0.05	<0.10		V0.037	<u> </u>	< 0.037
4,4'-DDE	<del> </del>						< 0.03	< 0.10				
4,4'-DDT	<c.091< td=""><td>&lt; 0.091</td><td>&lt; 0.091</td><td>&lt;0.091</td><td>&lt; 0.091</td><td>160:0&gt;</td><td>&lt; 0.05</td><td>&lt; 0.10</td><td></td><td>&lt; 0.091</td><td>10.001</td><td></td></c.091<>	< 0.091	< 0.091	<0.091	< 0.091	160:0>	< 0.05	< 0.10		< 0.091	10.001	
Dieldrin	<0.044	< 0.044	< 0.044	< 0.044	< 0.044	< 0.044	< 0.05	< 0.10		< 0.091	< 0.091	< 0.091
Endosulfan I	10.011	V0.044	\U.U++	10.044	₹0.044	<b>~0.044</b>	< 0.05	< 0.10		<0.044	< 0.044	< 0.044
Endosulfan II							< 0.05	< 0.20				
Enderulfan Sulfare							< 0.05	<0.10				
Endrin	< 0.039	< 0.039	< 0.039	< 0.039	- O 020	40.010						
Endrin Aldehyde	(0.03)	< 0.039	<b>~0.039</b>	(0.039	< 0.039	< 0.039	< 0.05	< 0.10	< 0.10	< 0.039	< 0.039	< 0.039
Heptachior	< 0.040	< 0.040	< 0.040	< 0.040	< 0.040	< 0.040	< 0.05	< 0.10				
Heptachlor Epoxide	< 0.032	< 0.032	< 0.040	< 0.632			< 0.05	< 9.10	< 0.10	<0.040	< 0.040	< 0.040
Methoxychlor	< 0.176	< 0.032	< 0.032	< 0.176	< 0.032	<0.032	< 0.05	< 0.10		< 0.032	< 0.032	< 0.032
Toxaphene	< 0.086	< 0.086			< 0.176	< 0.176	< 0.05	<1.0	<1.0	< 0.176	< 0.176	< 0.176
Aroclor-1016	< 0.050	< 0.050	<0.086 <0.050	< 0.036	< 0.086	< 0.086	<1.0	<2.0	<2.0	< 0.086	< 0.086	< 0.086
Aroclor-1221	< 0.054	< 0.054			< 0.050	< 0.050				< 0.050	< 0.050	< 0.050
Aroclor-1232	< 0.065	< 0.065	<0.054	< 0.054	< 0.054	< 0.051				< 0.054	< 0.054	< 0.054
Aroclor-1242	< 0.065			< 0.065	< 0.065	< 0.065				< 0.065	< 0.065	< 0.065
Aroclor-1248	< 0.000	< 0.065	< 0.065	< 0.065	< 0.065	< 0.065				< ∴065	< 0.065	< 0.065
Aroclor-1254	< 0.100	< 0.090	< 0.090	< 0.090	< 0.090	< 0.090				< 0.090	< 0.090	< 0.090
Aroclor-1260	< 0.100	<0.100 <0.100	< 0.100	<0.100	< 0.100	< 0.100				< 0.100	< 0.100	< 0.100
Arock 1268	< 0.100		<0.100 <0.100	<0.100	< 0.100	< 0.100				< 0.100	< 0.100	< 0.100
PCB-1016/1242	<u> </u>	<0.100	< 0.100	<0.100	< 0.100	< 0.100				< 0.100	< 0.100	< 0.100
PCB-1221	<del>  </del>						< 0.05	<1.0	<1.0			
PCB-1232	<del> </del>						< 0.03	<1.0	<1.0			
PCB-1248	<del> </del>						< 0.05	<1.0	<1.0			
PCB-1254	<del>  </del>						< 0.05	<1.0	<1.0			
PCB-1260	f						<.0.05	<1.0	<1.0			
PCB-1268	<del>  </del>						< 0.05	< 1.0	<1.0			
2,4-D	₹.0.29	< 0.29		< 0.29					<1.0			
Silvex(2,4,5-TP)	<0.34	< 0.34	<0.29 <0.34	< 0.29	< 0.29	< 0.29	< 2.0	<2.0	<2.0	< 0.29	< 0.29	< 0.29
Aldicarb	<0.50	< 0.34 < .50			< 0.34	<0.34	< 2.0	<.2.0	< 2.0	< 0.34	< 0.34	< 0.34
Carbofuran			<.50	< 0.50	<.50	<.50	< 10	< 10	< 10	< 0.50	<.50	<.50
1,2-Dibromo-3-Chloropropane	<5.0 <0.10	< 5.0	< 5.0	< 5.0	< 5.0	< 5.0	< 10	<10	< 10	<5.0	< 5.0	< 5.0
Ethylene dibromide	<0.10	< 0.10 < 0.10	<0.10	< 0.10	< 0.10	< 0.10				< 0.10	< 0.10	< 0.10
	₹0.10	< 0.10	< 0.10	< 0.10	< 0.10	< 0.10				<0.10	< 0.10	< 0.10
Trifluralia	<u></u>			l		l						

Parameter	MW2-D 7/26/99	MW2-D 9/9/99	MW2-D 11/1/99	MW3-D 1/26/99	MW3-D 9/9/99	MW3-D 11/1/99	MW4-D 7/26/99	MW4-D 9/9/99	MW4-D 11/1/99	MW5-D 2/9/99	MW5-D 11/1/99	MW6-D 7/26/99
Acenaphthene												
Acenaphthylene												
Alachior (lasso)	<1.0	< 1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	< 1.0
Atrazine	<1.0	<1.0	<1.0	< 1.0	<1.0	< : 0	<1.0	<1.0	<1.0	<1.0	<1.0	< 1.0
Anthracene												
Benzidine												
Benzo(a)anthracene												
Benzo(b)fluoramhene												
Benzo(k)fluoranthene												
Венго(в)ругеле												
Benzo(g,h,i)perylene												
Benzyl alcohol												
Benzyl butyl phthalate												
Bis(2-chluroethyl) ether												
Bis(2-chlorethoxy)methane												
Bis(2-ethylliexyf)phthalate	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	< 1.0
Bis(chloromethyl)ether	<1.0	<1.0			<1.0				<1.0	<1.0	<1.0	< 1.0
Bis(2-chlorolsopropyt)ether	\ \\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\	<u> </u>	\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\	<del></del>	~ 1.0	\ 21.0	71.0	``.0		\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\	\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\	~
4-Bromophenyl phenyl ether	<del> </del>		<b></b>		<del></del>							
4-Chloroaniline						<b> </b>						
	<del>                                     </del>				<del>                                     </del>							
2-Chloronaphthalene	<del> </del>		ļ	<b></b>								
4-Chlorophenylphenyl ether	<del> </del>				<del></del>							
Chrysenc	<del>                                     </del>		<b></b>	<u> </u>								
Dibenzo(a,b) anthracene	<del> </del>					ļ						
Di-n-butylphthalate	<1.0	< 1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.C	<1.0	<1.0	<1.0	<1.0
1,2-Dichlorobenzene	<b></b>			ļ <u></u>		<u> </u>						
1,3-Dichlorobenzene	<u> </u>					ļ						
1,4-Dichlorobenzene	ļ			<b></b>	ļ							
3,3-Dichlorobenzidine	ļ		ļ			<u> </u>			ļ			
Diethyl phthalate	<1.0	< 1.0	< 1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	< 1.0	<1.6
Dimethyl phthalate	<1.0	< 1.0	< 1.0	< 1.0	<1.0	< 1.0	<1.0	< 1.0	<1.0	<1.0	<1.0	<1.0
2,4-Dinitrotoluene	<u> </u>		ļ									
2,6-Dinitrotoluene	<u> </u>			<u> </u>								
Di-n-octylphthalate				·		<u> </u>		<u></u> .		<u> </u>		
Fluoranthene	<u> </u>				<u> </u>		L					
Fluorene	<u> </u>			1								
Hexachloro-1,3-butadiene												
Hexachlorocyclopentadiere												
Flexachloroethane												
Indenu(1,2,3 cd)pyrene	[						Ī					
Isophorone	<1.0	< 1.0	<1.0	< 1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0
2-Methylnaphthalene	<u> </u>	I	]	· · · · · ·	<del></del>	T	1	<u> </u>	T		<u> </u>	
Naphthalene	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0
2-Nitroaniline	T		<u> </u>	<u> </u>	<u> </u>	T	<u> </u>	<u> </u>	1	1	<u> </u>	-
3-Nitroaniline	<u> </u>	<u> </u>	<u> </u>	i		l		1	1	1	1	
4-titroaniline	T		T	]		1	<del>                                     </del>	<b>T</b>	l		<del> </del>	
Nitrobenzeile	1		1			1	1	1	<del>                                     </del>		<u> </u>	
N-Nitrosodimethylamine			1	<u> </u>	<u> </u>	1	<u> </u>	1	T	<del> </del>	<u> </u>	
N-Nitrosodiphenylamine	1		<del> </del>	<u> </u>	<del> </del>	t	<del> </del>	<del> </del>	<del> </del>	<del> </del>	<del>                                     </del>	
N-Nitrorodi-n-propylamine	<del> </del>		<del> </del>	<b> </b>	<del> </del>	<del></del>	<del> </del>		<del> </del>	<del> </del>	<u> </u>	<b></b>
Phenanthrene	<del>                                     </del>		<del>                                     </del>	l	<u> </u>	<del> </del>	<del> </del>	<del> </del>	<del>                                     </del>	<del> </del>	<del> </del>	<del></del>
Pyrene	†	<del></del>	<del> </del>	<del> </del>	<del></del>	<del> </del>	<del> </del>	<del> </del>	<del> </del>	<del> </del>	<del> </del>	<del></del>
Pyrkline	<del> </del>		<del> </del>	<del> </del>	<del> </del>	<del> </del>	<b> </b>	<del> </del>	<del> </del>	<del> </del>	<del> </del>	f
1,2,4-Trichlorobenzene	<del> </del>		<del> </del>	<del> </del>		<del> </del>		<del> </del>	<del> </del>	<del> </del>	<del> </del>	
	<del> </del>		<del> </del>	<del> </del>	<del> </del>	<del> </del>	<del> </del>	<del> </del>	<del> </del>	<del> </del>	<del> </del>	<b> </b>
Benzoic Acid	<del> </del> -		<del> </del>	<b> </b>		<del> </del> -	<del> </del> -	ļ	<del> </del>	<del> </del>	<del> </del>	<del> </del>
4-Chloro-3-methylphenol	<del> </del>		<del> </del>	<del> </del>	<del></del>	<del> </del> -	<del> </del> -	<del> </del>	<del> </del>	<del> </del>	<del> </del> -	<b> </b>
2-Chlorophenol	<del> </del>			ļ	<del> </del>	<b> </b>	<b></b>	ļ	ļ	<b> </b> -	<del> </del>	<b></b>
p-Cresol	<1.0	< 1.0	<1.0	< 1.0	<1.0	< 1.0	< 1.0	<1.0	< 1.0	<1.0	<1.0	<1.0
Cresols, Total	ļ		<b> </b>	<b></b>		ļ	ļ	<u> </u>	ļ	<b> </b>		<u> </u>
2,4-Dichlorophenol	ļ			<b></b>	<b></b>	<b></b>	<b></b>		<b> </b>	<u> </u>	<u> </u>	
2,4-Dimethylphenol		<b></b>	<u> </u>	ļ	ļ	<b> </b>	<b> </b>	<b></b>	<u> </u>	<b> </b>		
2,4-Dinitrophenol	<u> </u>	<u>L</u>	<u> </u>	<u>L</u>	<u></u>	<u> </u>	<u> </u>	<u>L</u>		<u></u>	1	
										-		

Parameter	MW2-D 7/26.99	MW2-D 9/9/99	MW2-D 11/1/99	MW3-D 7/26/99	MW3-D 9/9/99	MW3-D 11/1/99	MW4-D 7/26:99	MW4-D 9/9/99	MW4-D 11/1/99	MWS-D 9/9/99	MW5-D 11/1/79	MW6-D 7/26/99
2-Methyl-4,6-Ginitrophenol												•
2-Nicropteaol												
4-Nitrophenol												
Parathion	<1.0	<1.0	<1.0	< 1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0
Pentachlorophesol	<1.0	<1.0	<1.0	< 1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0
Phenol	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	71.0	<1.0	<1.0	<1.0
2,4,5-Trichloroptenol												
2,4,6-Trichlorophenol												
Aldrin	< 0.034	< 0.034	< 0.034	< 0.034	< 0.034	< 0.034	< 0.034	< 0.034	< 0.034	< 0.034	< 0.034	< 0.034
alpha-BHC												
benz-BHC												
gamma-BHC	< 0.025	< 0.025	< 0.025	< 0.025	< 0.025	< 0.025	< 0.025	< 0.025	< 0.025	< 0.025	< 0.025	< 0.025
delta-BHC												
Chlordane	< 0.037	< 6.037	< 0.037	< 0.037	< 0.037	< 0.037	< 0.037	< 0.037	< 0.037	< 0.037	< 0.037	< 0.037
4,4'-DDD	19.037	10.037	20.037	10.057	<u> </u>		10.05	40.037	30,037	10.03		
4,4'-DDE												
4.4'-PDT	< 0.091	< 0.091	< 0.091	< 0.091	< 0.091	<0.091	< 0.091	< 0.091	< 0.091	< 0.091	< 0.091	< 0.091
Dieldrin	< 0.044	< 0.044	< 0.044	< 0.044	< 0.044	<0.044	< 6.044	< 0.044	< 0.044	< 0.044	< 0.044	< 0.044
Endesulfan I			- 30.017		10.044	1,0,014	10.011		70.04	10.04	70.04	19:011
Endosulfan II												
Endosulfan Sulfate	<b> </b>	<b></b>										
Endrin	< 0.039	< 0.039	< 0.039	< 0.039	< 0.039	< 0.035	< 0.039	< 0.019	< 0.039	< 0.039	< 0.039	<0.039
Endria Aldehyde	\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\	C 0.039	(0.039	C0.033	< 0.039	< 0.035	\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\	< 0.039	- (0 039	(0.039	(0.039	4.10,000
Heptachlor	< 0.040	< 0.040	<0.040	< 0.040	< 0.040	< 0.040	< 0.040	< 0.040	< 0.040	< 0.040	<0.040	< 0.040
	< 0.032	< 0.032	< 0.032	< 0.032	< 0.032	< 0.032	< 0.032	< 0.032	< 0.040	< 0.032	< 0.032	< 0.032
Heptachlor Epoxide	<0.176	< 0.176	< 0.176	< 0.176			< 0.176		< 0.176			< 0.176
Methoxychlor	< 0.036				< 0.176	< 0.176		< 0.176		< 0.176	< 0.176	
Totaphene Aroclor-1016	< 0.050	< 0.086 < 0.060	< 0.086 < 0.080	< 0.01/3	< 0 386	< 0.050	< 0.086 < 0.050	< 0.086	< 0.086 < 0.050	< 0.086 < 0.050	< 0.086	< 0.086 < 0.050
	< 0.054			< 0.054	1.050</td <td></td> <td></td> <td></td> <td></td> <td></td> <td></td> <td></td>							
Arrickir-1221 Arcs for -1232	< 0.054	< 0.054	< 0.054	< 0.065	< 0.054	<0.054	< 0.054	< 0.054	< 0.054	< 0.054	< 0.034	< 0.054
Arcclor-1242	< 0.065		< 0.065		< 0.065	< 0.065	< 0.065	< 0.065	< 0.063	< 0.065	< 0.065	< 0.065
		< 0.065	< 0.065	< 0.055	< 0.065	< 0.065	< 0.065	< 0.065	< 0.065	< 0.065	< 0.065	< 0.063
Arccior-1 1/18	< 0.090	< 0.090	< 0.090	< 0.090	< 6.090	< 0.090	< 0.090	< 0.090	< 0.090	< 0.090	< 0.090	< 0.090
Aroclor-1254	< 0.100	< 0.100	< 0.100	< 0.100	< 0.100	< 0.100	< 0.100	< 0.100	< 0.100	< 0.100	< 0.100	< 0.100
Arc:lor-1260	< 0.100	< 0.100	< 0.100	< 0.100	< 0.100	< 0.100	< 0.100	< 0.100	< 0.100	< 0.100	< 0.100	< 0 100
Aroclor-1268	< 0.100	< 0.100	< 0.100	< 0.100	< 0.100	< 0.100	< 0.100	< 0.100	< 0.100	< 0.100	< 0.100	< 0.100
PCB-1016/1243		<del> </del>	<del> </del> -	<u> </u>	<b> </b>		<b></b>		ļ		·	
PCB-1221	ļ		<b></b>		ļ		<b> </b>		ļ			
PCB-1232	<del> </del>	<u> </u>	<del></del> -				}					
PCD-1248	<b></b>	ļ			<del> </del>	<u> </u>	<b> </b>	ļ	ļ			
PCB-1254	<del> </del>		<u>;</u>				ļ	<del></del>	<b></b>			
PCB-1260	ļ	<b> </b>					ļ					
PCB-1268			ļ		<u> </u>	ļ.———	ļ		ļ		<b></b>	
<u>{2,4-D</u>	< 0.29	<del></del>		< 0.29	< 0.29	< 0.29	< 0.29	<.0.29		< 0.29	< 0.29	< 0.29
Silves (2,4,5-TP)	< 0.34	< 0.34		< 0.34	< 0.34	< 0.34	< 0.34	< 0.34	†-·	< 0.34	< 0.34	< 0.34
Aldicarb	< 0.50	< .50	<del> </del>	< 0.50	< .50	< .50	< 0.50	< .50			< .50	<0.50
Carboluran	≤5.0	< 5.0	<del></del>	< 5.0	< 5.0	< 5.0	< 5.0	< 5.0			< 5.0	< 5.0
1 2 Dibromo 3 Chloropropure	< 0.10	< 0.10	f	< 0.10	< 0.10	< 0.10	< 0.10	< 0.10	< 0.10	< 0.10	< 0.10	< 0.10
Ethylene dibromide	< 0.10	< 3.10	< 0.10	< 0.10	< 0.10	< 0.10	< 0.10	< 0.10	< 0.10	<0.i0	< 0.10	< 0.10
Triffuralin	<u></u>	<u> </u>	L	<u> </u>	L	<u> </u>	1	l	<u></u>	<u></u>	<u> </u>	

Parameter	MW6-D 9/9/99	MW6-D 11/1/99	MW7-D 7/26/99	MW7-D 9.9/99	MW7-D 11/1/99	MW8-D 7/26/99	MW8-D 9/9/99	MW8-D 11/1/99	MW9-D 7/26/99	MW9-D 9/9/99	MW9-D 11/1/99	MW1-W5 7/26/99
Acenephthese					·							l
Aceraphthylene												
Alachlor (lasso)	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	< 1.0	<1.0	<1.0	<1.0
Atrazine	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	< 1.0	<1.0	<1.0	<1.0	<1.0	
Anthracene												
Benzidine												
Benzo(a)anthracene												
Benzo(b)fluoranthene												
Benzo(k)fluoranthene										7. 7. 7. 7.		
Benzo(a)pyrene												
Benzo(3, h, i)perylere												
Benzyi alcohol												
Benzyl butyl phthalate												
Bis(2-chloroethyl) ether												
Bis(2-chlorethoxy)mediane		<b>6</b> 49										
Bis(2-ethylhexyl)phthalate	<1.0	< 1.0	< 1.0	< 1.0	<1.0	<1.0	<1.0	<1.0	<1.0	< 1.0	< 1.0	<1.0
Bis(chloromethyf)ether	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	< 1.0	<1.0	<1.0
Bis(2-chlorohopropyf)ether				17.0		****	~1.0	<u></u>		<u></u>	<u> </u>	<u> </u>
4-Bromophenyl phenyl ether												<del></del>
4-Chloroaniline	1											<del></del>
2-Chloronaphthalene	-											
4-Chlorophenylphenyl ether	1											
Chrysene	1											
Dibenzo(a,h) anthracene	1											
Di-n-butylphthalate	<1.0											
1,2-Dichlorobenzene	1 21.0	<1.0	< 1.0	:1.0	<1.0	< 1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0
1,3-Dichlorobeniene	+											
~ <del>```</del>	<del></del>											<del></del>
1,4-Dichlorobenzene											-	
3,3-Dichlorobeazidine	-											·
Directly obstalate	<1.0	<1.0	<1.0	< 1.0	<1.0	< 1.0	< 1.0	< 1.0	<1.0	<1.0	< 1.0	< 1.6
Dimethyl phthalate	<10	<1.0	<1.0	<1.0	<1.0	< 1.0	< 1.0	< 1.0	< 1.0	< 1.0	< 1.0	<1.0
2,4-Dinkrotoluene	-											
2,6-Dinitrosolvene						<del>.</del> -						
Di-n-octylphthalate Fluoranthene	<del> </del>											
Fluorene					[							
······································	<del>  </del>											
Hexachloro-1,3-butadiene	+											
Hexachlorocyclopentadiene  Manachlorocyclopentadiene												
Hexachloroethane Indeno(1,2,3-cd)pyrene	<del>  </del>											
Isophorone	1											
2-Methylnaphthalene	<1.0	< 1.0	<1.0	<1.0	<1.0	< 1.0	< 1.0	<1.0	<1.0	<1.0	< 1.0	<1.0
Naphibalene 7-Nitroaniline	<1.0	< 1.0	<1.0	< 1.0	<1.0	< 1.0	<1.0	<1.0	<1.0	<1.0	<1.0	< 1.0
3-Naroanilina	<del> </del>								1			
	<del></del>											
4-Nitroaniline	<del> </del>				.,							
Niuoberzene												
N-Nitrosodimed ylamine										l		
N-Nitrosodiphenylamine												
N-Nitrosodi-n-propylamine								l				
Phenanthrene	1				I	7		1				
Pytere	1		l	I		1						
Pyridine	1		]						1			
1.2.4-Trichlorobenzene	<u> </u>											
Henzoic Acid	T											
4-Chloro-3-methylpherial	T											
2-Chlorophenol	T											
p-Cresol	<10	<1.0	<1.0	<1.0	<1.0	<1.0						
Cresols, Total	┟──┴┼				3	~ 1.0	< 1.0	< 1.0	< 1.0	<1.0	< 1.0	<1.0
2,4-Dichlorophenol	1											
	<del>  -</del>											
2,4 Direchylphenol												

Parameter '	MW6-D 9/9/99	MW6-D 11/1/99	MW7-D 7/26/99	MW7-D 9/9/99	MW7-D 11/1/99	MW8-D 7/26/99	MW8-D 9/9/99	MW8-D 11/1/99	MW9-D 7/26/99	MW9-D 9/9/99	MW9-D 11/1/99	MW1-WT 7/26/99
2-Methyl-4,6-dinitrophenol												
2-Nitrophenol												
4-Nitrophenol												
Parathion	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	< 1.0	<1.0	<1.0	<1.0	<1.0
Pentachlorophenol	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	< 1.0
Pherest	<1.0	< 1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	< 1.0	<1.0	<1.0	< 1.0
2,4,5-Tricklorophenol												
2,4,6-Trichlorophenol												
Aldrin	< 0.034	< 0.034	< 0.034	< 0.034	< 0.034	< 0.034	< 0.034	< 0.034	< 0.034	< 0.034	< 0.034	< 0.034
alpha-BHC												
beta-BHC												
датта-ВНС	< 0.025	< 0.025	< 0.025	< 0.025	< 0.025	< 0.025	< 0.025	< 0.025	< 0.025	< 0.025	< 0.025	< 0.025
delta-BHC												
Chlordane	< 0.037	< 0.037	< 0.037	< 0.037	< 0.037	< 0.037	< 0.037	< 0.037	< 0.037	< 0.037	< 0.037	< 0.037
4,4'-DDD												
4,4'-DDE												
4,4'-DDT	< 0.091	< 0.091	< 0.091	< 0.091	< 0.091	< 0.091	< 0.091	< 0.091	< 0.091	< 0.091	< 0.091	< 0.091
Dieldrin	< 0.044	< 0.044	< 0.044	< 0.044	< 0.044	< 0.044	< 0.044	< 0.044	< 0.044	< 0.044	< 0.044	< 0.044
Endosulfan I												
Endosulfan II												
Endosulfan Sulfate												
Endrin	< 0.039	< 0.039	< 0.039	< 0.039	< 0.039	< 0.039	< 0.039	< 0.039	< 0.039	< 0.039	< 0.039	< 0.039
Endrin Aldehyde												
Heptachlor	< 0.040	< 0.040	< 0.040	< 0.040	< 0.010	< 0.040	< 0.040	< 0.040	< 0.040	< 0.040	< 0.040	< 0.040
Heptachic. Epoxide	< 0.032	< 0.032	< 0.032	< 0.032	< 0.032	< 0.032	< 0.032	< 0.032	< 0.032	< 0.032	< 0.032	< 0.032
Methoxychlor	< 0.176	< 0.176	< 0.176	< 0.176	< 0.176	< 0.176	< 0.176	< 0.176	< 0.176	< 0.176	< 0.176	< 0.176
Toxaphene	< 0.086	< 0.086	< 9.086	< 0.086	< 0.086	< 0.086	< 0.086	< 0.086	< 0.086	< 0.086	< 0.086	< 0.086
Aroclor-1016	< 0.050	< 0.050	< 0.050	< 0.050	< 0.050	< 0.050	< 0.050	< 0.050	< 0.050	< 0.050	< 0.050	< 0.050
Aroclor-1221	< 0.054	< 0.054	< 0.054	< 0.054	< 0.054	< 0.054	< 0.054	< 0.054	< 0.054	< 0.054	< 0.054	< 0.054
Aroclor-1232	< 0.065	< 0.063	< 0.065	< 0.065	< 0.065	< 0.065	< 0.065	< 0.065	< 0.065	< 0.065	< 0.065	< 0.065
Aroclor-1242	< 0.063	< 0.065	< 0.065	< 0.065	< 0.065	< 0.065	< 0.065	< 0.065	< 0.065	< 0.065	< 0.065	< 0.065
Aroclor-1248	< 0.090	< 0.090	< 0.090	< 0.090	< 0.090	< 0.090	< 0.090	< 0.090	< 0.090	< 0.090	< 0.090	< 0.090
Arocler 1254	< 0.100	< 0.100	< 0.100	< 0.100	< 0.100	< 0.100	< 0.100	< 0.100	< 0.100	< 0.100	< 0.100	< 0.100
Aroclur-1260	< 0.100	< 0.100	< 0.100	< 0.100	< 0.100	< 0.100	< 0.100	< 0.100	< 0.100	< 0.100	< 0.100	< 0.100
Aroclor-1268	< 0.100	< 0.100	< 0.100	< 0.100	< 0.100	< 0.100	< 0.100	< 0.100	< 0.100	< 0.100	< 0.100	< 0.100
PCB-1016/1242												,
PCB-1221												
PCB-1232												
PCB-1248												
PCB-1254												
PCB-1260												
PCB-1268												
2,4·D	< 0.29	< 0.29	< 0.29	< 0.29	< 0.29	< 0.29	< 0.29	< 0.29	< 0.29	< 0.29	< 0.29	< 0.29
Silvex(2,4,5-TP)	< 0.34	< 0.34	< 0.34	< 0.34	< 0.34	< 0.34	< 0.34	< 0.34	< 0.34	< 0.34	< 0.34	< 0.34
Aldicarb	< .50	< .50	< 0.50	< .50	< .50	< 0.50	< .50	< .50	< 0.50	<.50	<.50	< 0.50
Carbofuran	< 5.0	< 5.0	< 5.0	< 5.0	< 5.0	< 5.0	< 5.0	< 5.0	< 5.0	<5.0	< 5.0	<5.0
1,2-Dibromo-3-Chloropropane	< 0.10	< 0.10	< 0.10	< 0.10	< 0.10	< 0.10	< 0.10	< 0.10	< 0.10	< 0.10	<0.10	< 0.10
Ethylene dibromide	< 0.10	< 0.10	< 0.10	< 0.10	< 0.10	< 0.10	< 0.10	<0.10	< 0.10	< 0.10	< 0.10	< 0.10
Trifluralio	1											

Parameter	MW1-WT 9/9/99	MW1-WT 11/1/99	MW2-WT 3/2/00	MW2-WT 7/26/99	MW2-WT 9/9/99	MW2-WT 11/1/99	MW3-WT 7/26/99	MW3-WT 9/9/99	MW3-WT 11/1/99	MW4-WT 7/26/99	MW4-WT 9/9/99	MW4-WT
Acenaphthene		******	3/2/00	.,20,22	213132	11/1/55	1,2433	2,2,2,2	11/1/2	1120.77		22.25
Acenaphthylene												····
· · · · · · · · · · · · · · · · · · ·					41.0					<1.0	<1.0	-110
Alachior (lasso)	<1.0	<1.0	<1.0	<1.0	<1.0		<1.0	<1.0	<1.0			<1.0
Atrazine	<1.0	<1.0	< 1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	< 1.0
Anthracene Benzidine											·	
Benzo(a)anthracene												
Benzo(b)fluorantiene			<u> </u>									
Benzo(k)fluoranthens												
Benzo(a)pyrene												
Benzo(g,h,i)perylene										ļ		
Benzyl alcohol						ļ						
Benzyl butyl phthalate			<u> </u>			<u> </u>						
Bis(2-chloroethyl) ether												
Bis(2-chlorethoxy)methane								L				
Bis(2-ethylhexyl)phthalate	<1.0	<1.0	< 1.0	< 1.0	<1.0	< 1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0
Bis(chloromethyl)ether	<1.0	<1.0	< 1.0	<1.0	<1.0	<1.0	< 1.0	<1.0	<1.0	<1.0	<1.0	<1.0
Bis(2-chloroisopropyl)ether		ļ	<b></b>	ļ		ļ			ļ	ļ		
4-Bromophenyi phenyi ether			<u></u>		ļ	<u> </u>	<u> </u>	<u> </u>		<b></b>		<u> </u>
4-Chloroaniline	_		<u> </u>		<b></b>		<b></b>					
2-Chloronaphthalene												
4-Chlorophenylphenyl ether			L		<u> </u>							
Chrysene												
Dibenzo(a,h) anthracene												
Di-n-buty!phthalate	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0
1,2-Dichlorobenzene												
1,3-Dichlorobenzene							[		T			
1,4-Dichlorobenzene										l		
3,3-Dichlorobenzidine												
Diethyl phthalate	<1.0	<1.0	<1.0	<1.0	<1.0	< 1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0
Dimethyl phthalate	<1.0			<1.0			1					<1.0
2,4-Dinitrotoluene				<u> </u>					<u></u>	<u></u>		<u>``-</u>
2,6-Dinitrotoluene			<u> </u>			l			i		<del> </del>	<b></b>
Di-n-octylphthalate			<del> </del>					<b></b>		ļ	<u> </u>	
Fluoranthene						<del> </del>				·	·	<b></b>
Fluorene		<del></del>	<u> </u>	<b></b>		<del> </del>	<del> </del>			<del> </del>	<del></del>	<del></del> -
Hexachloro-1,3-butadiene		<b></b>		<del> </del>	<b></b>	† <del></del> -	<del> </del>		<del> </del> -		<del> </del>	<del> </del>
Hexachlorocyclopentadiene			<del></del>		<del></del>	<del></del>	<del>                                     </del>	····	<del> </del>			<del></del>
Hexachleroethane		<del></del>		<del> </del>		<del> </del>	<del> </del>	<del> </del>		<del> </del>	<del> </del>	
Indeno(1,2,3-cd)pyrene		<b></b>		<b></b>			<del> </del>			<del> </del>	<del> </del>	<del></del>
Isophorone	<1.0	<1.0	< 1.0	< 1.0	< 1.0	<1.0	<1.0	<1.0	<1.0	<1.0	-10	
2-Methylnaphthalene		<u></u>	1	<u> </u>	\ \ \ \ \ \ \ \ \ \ \ \ \ \ \ \ \ \ \		<u> </u>	<u> </u>	<u> </u>	1 1.0	< 1.0	<1.0
			<10	<1.0	~10		<1.0					
Naphthalene 2-Nitroaniline	< 1.0	<1.0	<1.0	<del>`</del> '	<1.0	<1.0		<1.0	<1.0	<1.0	< 1.0	< 1.0
3-Nitroaniline			<del> </del>	<del> </del>	<del> </del>	<del> </del>	<del> </del>	<del> </del>	<del></del> -	<del> </del>	<del> </del>	<del> </del>
		<del> </del>	<del> </del>	<del> </del>	<del> </del>	<del> </del>	<del> </del>	<del> </del>	<del> </del>	<del> </del>	<del> </del>	<del> </del>
4-Nitroaniline		ļ	<del> </del> -	<del> </del>	<del> </del>	<del> </del>	<del> </del>	<del> </del>	<del> </del>	<del> </del> -	<del> </del>	<del> </del>
Nitrobenzene			<del> </del>	<del></del>	<del> </del>	ļ	<del> </del>	<del></del>	<del> </del>	<del> </del>	<del> </del>	<del> </del>
N-Nitrosodimethylamine			<del> </del>	<del> </del>	<del> </del>	<del> </del>	<del> </del> -	<del> </del>	ļ	<del> </del>	<b></b>	<del> </del>
N-Nitrosodiphenylamine	<del></del>	<del></del>	<del> </del>	<del> </del>	ļ	<del> </del>		<del> </del>	<del> </del>	<del> </del>	ļ	ļ
N-Nitrosodi-n-propylamine		<b></b>	<del> </del>		<del> </del> -	<del> </del>	<del> </del>	<b> </b>	<b> </b>	<u> </u>	<del> </del>	<del> </del>
Phenanthrene			<del> </del>		<b> </b>	<b></b>	<del> </del> -		<del> </del>	<del> </del>	ļ	<b></b>
Pyrene		ļ	<del> </del>	<del> </del>	ļ	<del> </del>	ļ	<del> </del>	<del> </del>	ļ	<b> </b> -	<u> </u>
Pyridine		<del></del>	<b> </b>	<del> </del>	ļ	<b></b>	ļ	<del> </del>	ļ	<b> </b>		ļ
1,2,4-Trichlorobenzene			ļ	<del> </del>	<b> </b>	ļ	ļ	<b> </b>		<b> </b> -	ļ	ļ
Benzoic Acid			ļ	<b> </b>	ļ	ļ	ļ	<b> </b> _	ļ	<b> </b>	<u> </u>	<u> </u>
4-Chloro-3-methylphenul			<b></b>	<b></b>	<b> </b>	<u> </u>	<u> </u>	<u> </u>				L
2-Chlorophenol				<u> </u>			ļ		<u></u>			
p-Cresol	< 1.0	<1.0	<1.0	< 1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.
Cresols, Total				<u> </u>							T	1
2,4-Dichlorophenol									I		<u> </u>	
2,4-Dimethylphenol			I	l	<u> </u>	1	I	1		T	1	<del>                                     </del>
2,4-Dinitrophenol				<del></del>	·	<del></del>	<del></del>	1	<del> </del>	<del> </del>	<del> </del>	<del> </del>

Parameter	MW1-WT	MW1-WT	MW1-WT	MW2-WT		MW2-WT	AIW3-WT			MW4-WT	MW4-WT	MW4-WT
	9/9/99	11/1/99	3/2/00	7/25/99	9/9/99	11/1/99	7/26/99	9/9/99	11/1/99	7/26/99	9/9/99	11/1/99
2-Methyl-4,6-dinitrophenol												
2-Nitrophenol												ļ
4-Nitrophenol												
Parathion	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0
Pentachlorophenol Phenol	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	< 1.0	<1.0	<1.0	<1.0	<1.0	< 1.0
	<1.0	<1.0	<1.0	< 1.0	<1.0	<1.0	<1.0	.0</td <td>&lt;1.0</td> <td>&lt;1.0</td> <td>&lt;1.0</td> <td>&lt; 1.0</td>	<1.0	<1.0	<1.0	< 1.0
2,4,5-Trichlorophenol		<u> </u>			<del></del>							
2,4,6-Trichlorophenol					<u> </u>							
Aldrin	< 0.034	< 0.034	< 0.034	< 0.034	< 0.034	< 0.034	< 0.034	< 0.034	< 0.034	< 0.034	< 0.034	< 0.034
alpha-BHC						ļ						
beta-BHC					<del> </del>							ļ
gamma-BHC	< 0.025	< 0.025	< 0.025	< 0.025	< 0.025	< 0.025	< 0.025	< 0.025	< 0.025	< 0.025	< 0.025	< 0.035
delta-BHC						<b></b>			i			
Chlordane	<0.037	< 0.037	< 0.037	< 0.037	< 0.037	< 0.037	< 0.037	< 0.037	< 0.037	< 0.037	< 0.037	< 0.037
4,4'-DDD						<b></b>						
4,4'-DDE				ļ								
4,4'-DDT	< 0.09!	< 0.091	< 0.091	< 0.091	< 0.091	< 0.091	< 0.091	< 0.091	< 0.091	< 0.091	< 0.091	< 0.091
Dieldrin	< 0.044	< 0.044	< 0.044	< 0.044	< 0.044	< 0.044	< 0.044	< 0.044	< 0.044	< 0.044	< 0.044	< 0.044
Endosulfan ?					ļ							
Endosulfan II		ļ				<b></b>						
Endosulfan Sulfate		ļ										
Endrin	< 0.039	< 0.039	< 0.039	< 0.039	< 0.039	< 0.039	< 0.039	< 0.039	< 0.039	< 0.039	< 0.039	< 0.039
Endrin Aldehyde		ļ <u>.</u>				ļ						
Heptachlor	< 0.040	< 0.040	< 0.340	< 0.040	< 0.040	< 0.040	< 0.040	< 0.040	< 0.040	< 0.040	< 0.040	< 0.040
Heptachlor Epoxide	< 0.032	< 0.032	< 0.032	< 0.032	< 0.032	< 0.032	< 0.032	< 0.032	< 0.032	< 0.032	< 0.032	< 0.032
Methoxychlor	< 0.176	< 0.176	< 0.176	< 0.176	< 0.176	< 0.176	<0.176	< 0.176	< 0.176	< 0.176	< 0.176	< 0.176
Toxaphene	< 0.086	< 0.086	< 0.086	< 0.086	< 0.086	< 0.086	< 0.086	< 0.086	< 0.086	< 0.086	< 0.086	< 0.086
Aroclor-1016	< 0.050	< 0.050	< 0.050	< 0.050	< 0.050	< 0.050	< 0.050	< 0.050	< 0.050	< 0.050	< 0.050	< 0.050
Aroclor-1221	< 0.054	< 0.054	< 0.054	< 0.054	< 0.054	< 0.054	< 0.054	< 0.054	< 0.054	< 0.054	< 0.054	< 0.054
Aroclor-1232	< 0.065	< 0.065	< 0.065	< 0.065	< 0.065	< 0.065	< 0.065	< 0.065	< 0.065	< 0.065	< 0.055	< 0.065
Aroclor-1242	< 0.065	< 0.065	< 0.065	< 0.065	< 0.065	< 0.065	< 0.065	< 0.065	< 0.065	< 0.065	< 0.065	< 0.055
Aroclor-1248	< 0.090	< 0.090	< 0.090	< 0.090	< 0.090	< 0.090	< 0.090	< 0.090	< 0.090	<.0.690	< 0.090	< 0.090
Aroclor-1254	<0.100	< 0.100	2.100	< 0.100	< 0.100	< 0.100	< 0.100	< 0.100	< 0.100	< 0.100	< 0.100	< 0.100
Aroclor-1260	< 0.100	< 0.100	< 0.100	< 0.100	< 0.100	< 0.100	< 0.100	< 0.100	< 0.100	< 0.100	< 0.100	< 0.100
Aroclor-1268	< 0.100	< 0.100	< 0.100	< 0.100	< 0.100	< 0.100	< 0,100	< 0.100	< 0.100	< 0.100	< 0.100	< 0.100
PCB-1016/1242		ļ		<del></del> -	<del> </del>							
PCB-1221		ļ		ļ								ļ
PCB-1232		· <del></del>			<del> </del>				ļ			ļ
PCB-1248		}	<b> </b>	<u></u>	}- <u></u> -				ļ			ļ
PCB-1254					<b> </b> -							ļ
PCB-1260				<b></b>	<del> </del>							ļ
PCB268			-0.50		<del></del>			<b></b>	<b></b> _			
2,4-D	< 0.29	< 0.29	< 0.29	< 0.29	< 0.29	< 0.29	·:0.29	< 0.29	< 0.29	< 0.29	< 0.29	< 0.29
Silvex(2,4,5-TP)	<0.34	< 0.34	< 0.34	< 0.34	< 0.34	< 0.34	< 0.34	< 0.34	< 0.34	< 0.34	< 0.34	< 0.34
Aldicarb	<.50	<.50	<.50	< 0.50	< .50	<.50	< 0.50	< .50		< 0.50	<.50	<.50
Carbofuran	< 5.0	< 5.0	< 5.0	< 5.0	< 5.0	< 5.0	< 5.0	< 5.0	< 5.0	<5.0	<5.0	< 5.0
1,2-Dibromo-3-Chloropropane	< 0.10	< 0.10	< 0.10	< 0.10	< 0.10	< 0.10	< 0.10	< 0.10		<0.10	< 0.10	< 0.10
Ethylene dibromide	< 0.10	< 0.10	<0.10	< 0.10	< 0.10	< 0.10	< 0.10	< 0.10	<0.10	< 0.10	< 0.10	< 0.10
Trifluralin			L	L	l	L	L	L	l		L	

Parameter	MW9-WT 7/26/99	MW9-WT 9/9/99	MW9-WT 11/1/99	LW-1 7/24/98	LW-1 10/26/98	LW-1 7/26/99	LW-1 9/9/99	LW-1 11/1/99	LW-3 7/26/99	LW-3 9/9/99	LW-3 11/1/99	SPRING 7/26/99
Acenaphthene	1			< 20	< 10							
Acenaphthylene				<20	< 10						,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,	
Alachior (lasso)	<1.0	<1.0	<1.0	< 0.5	< 0.5	<1.0	<1.0	<1.0	<1.0	<1.0	< 1.0	<1.0
Atrazine	<1.0	<1.0	<1.0	<0.5	< 0.5	<1.0	<1.0	< 1.0	<1.0	< 1.0	< 1.0	<1.0
Anthracene				<20	< 10							
Benzidine				<200	<100							
Benzo(a)anthracene				< 20	< 10							
Benzo(b)fluoranthene				< 20	< 10							
Benzo(k)fluoranthene				< 20	< 10							
Benzo(a)pyrene				< 20	< 10							
Benzo(g,h,i)perylene				< 20	< 10							
Benzyl alcohol		<u> </u>		<20	< 10							
Benzyl butyl phthalate		<u> </u>		< 20	< 10							
Bis(2-chloroethyl) ether	_	<b> </b>		< 20	< 10							
Bis(2-chlorethoxy)methane	<u> </u>			< 20	< 10							
Bis(2-ethylhexyl)phthalate	<1.0	<1.0		< 20	< 10	<1.0	<1.0	<1.0	<1.0	<1.0	< 1.0	<1.0
Bis(chloromethyl)ether	<1.0	<1.0	<1.0			<1.0	<1.0	< 1.0	<1.0	< 1.0	<1.0	<1.0
Bis (2-chloroisopropyl)ether	<del> </del>	ļ	<b> </b>	< 20	< 10							
4-Bromophenyl phenyl ether		<del> </del>	<b> </b>	< 20	< 10							
4-Chloroaniline	<del>- </del>	ļ <u>.</u>	<del> </del>	< 20	< 10	<b> </b>						
2-Chloronaphthalene			ļi	< 20	< 10	<del> </del>						
4-Chlorophenylphenyl ether		<del> </del>	<b></b>	<20	< 10		<del></del>			ļi		
Chrysene		<del> </del> -		< 20	< i0							
Dibenzo(a,h) anthracene Di-n-butylphthalate				< 20	< 10							
1,2-Dichlorobenzene	<1.0	<1.0	< 1.0	< 20	< 10		<1.0	< 1.0	<1.0	< 1.0	< 1.0	<1.0
1,3-Dichlorobenzene		<del> </del>		<20 <20	< 10 < 10							
1,4-Dichlorobenzene	<del> </del>	<del> </del>	ļ	<20	< 10					<del> </del>		
3,3-Dichlorobenzidine	<del></del>	<del>                                     </del>	<del> </del>	< 100	<50							
Diethyl phthalate	<1.0	<1.0	<1.0	<20	< 19	<1.0	< 1.0	< 1.0	<1.0	<1.0	< 1.0	< 1.0
Dimethyl phthalate	<1.0			< 20	< 10		<1.0	<1.0	<1.0	<1.0	<1.0	<1.0
2,4-Dinitrotoluene	1	<del>```</del>	1	< 20				<u> </u>		<u> </u>	<u></u>	<u> </u>
2,6-Dinitrotoluene				<20	< 10							
Di-n-octylphthalate				< 20	<10							
Fluoranthene				<20	< 10							
Fluorene				< 20	< 10							
Hexachloro-1,3-butadiene				< 20	< 10	1					·	
Hexachiorocyclopentadiene				< 40	<20							
Hexachloroethane				< 20	< 10							
Indeno(1,2,3-cd)pyrene				< 20	< 10							
Isophorone	<1.0	< 1.0	<1.0	< 20	< 10	<1.0	<1.0	<1.0	< 1.0	<1.0	<1.0	<1.0
2-Methylnaphthalene				< 2.0	< 10							
Naphihalene	<1.0	<1.0	< 1.0	< 20	< 10	<1.0	< 1.0	< 1.0	< 1.0	< 1.0	<1.0	< 1.0
2-Nitroanlline			<u> </u>	< 20	< 10							
3-Nitroaniline		<b> </b>		< 20	< 10							
4-Nitroaniline		<u> </u>	<b> </b>	< 20	< 10						<u> </u>	
Nitrobenzene			ļi	< 20	< 10							
N-Nitrosodimethylamine				< 20	< 10							
N-Nitrosodiphenylamine	_	ļ		< 20	< 10					<u> </u>		
N-Nitrosodi-n-propylamine				< 20	< 10			<u> </u>				
Phenanthrene		<b> </b>	<b></b>	< 20	< 10				ļ	<b> </b>		<u> </u>
Pyrene		<b>}</b>		< 20	< 10	<del></del>				ļ	ļ	ļ
Pyridine	<del></del>	<b> </b>	<b> </b> _	< 20	< 10				<b></b>	<u> </u>	<b> </b> .	
1,2,4-Trichlorobenzene	<del></del>	ļ	ļ	< 20	< 10		<b></b>		<u> </u>	<u> </u>	ļ	
Benzoic Acid		<b> </b>	<b> </b>	<40	< 20		ļ		<b> </b>		ļ	
4-Chloro-3-methylphenol	<del></del>	<b> </b>		< 20		·			ļ	ļ	<u> </u>	
2-Chlorophenoi	_		ļ	< 20	< 10				ļ. <u>.                                   </u>	<u> </u>		
p-Cresol	< 1.0	< 1.0	< 1.0			<1.0	<1.0	<1.0	<1.0	<1.0	< 1.0	<1.0
Cresols, Total		<b> </b>	ļ	< 20	< 10	ļ			<u> </u>			
2,4-Dichlorophenoi		<b> </b>	<u> </u>	< 20					<u> </u>	<u> </u>		
2,4-Dimethylphenol	<b></b>	ļ	<b>}</b>	< 20						ļ	<u> </u>	<u> </u>
2,4-Dinitrophenol		l	L	< 20	< 10	1		L	L	<u></u>	L	

Parameter	MW9-WT 7/26/99	MW9-WT 9/9/99	MW9-WT 11/1/99	LW-1 7/24/98	LW-1 10/26/98	LW-1 7/26/59	LW-1 9/9/99	LW-1 11/1/99	LW-3 7/26/99	LW-3 9/9/99	LW-3 11/1/99	SPRING 7/26/99
2-Methyl-4,6-dinitrophenol				< 20	< 10							
2-Nitrophenol				<20	<10							
4-Nitrophenol				<20	<10							
Parathion	<1.0	< 1.0	<1.0	< 0.5	< 0.5	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0
Pentachlorophenol	<1.0	<1.0	<1.0	<20	< 10	<1.0	<1.0	<1.0	<1.0	< 1.0	<1.0	<1.0
Phenol	<1.0	<1.0	<1.0	<20	<10	< 1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0
2,4,5-Trichlorophenol				<20	< 10							
2,4,6-Trichlorophenol				<20	< 10							
Aldrin	< 0.034	< 0.034	< 0.034	<0.10	< 0.10	< 0.034	< 0.034	< 0.034	< 0.034	< 0.034	<0.034	< 0.034
alpha-BHC				<0.10	< 0.10	10,000	10.051	10.031	- 44.05.		10.051	10.001
beta-BHC				< 0.10	< 0.10							
gan.ma-BHC	< 0.025	< 0.025	< 0.025	< 0.10	< 0.10	< 0.025	< 0.025	< 0.025	< 0.025	< 0.025	< 0.025	< 0.025
delta-BHC				<0.10	< 0.10		V 0.023	V0.025	10.025	<u> </u>	1,0,023	
Chlordane	< 0.037	< 0.037	< 0.037	< 0.4	<1.0	< 0.037	< 0.037	< 0.037	< 0.037	< 0.037	< 0.037	<0.037
4,4'-DDD		1 1	10.037	<0.10	< 0.10	<b>40.937</b>	<b>VU.037</b>	<b>\\0.037</b>	V0.037	<u> </u>	< 0.037	20.037
4,4'-DDE				<0.10	<0.10							
4,4'-DDT	< 0.091	< 0.091	< 0.091	<0.10	<0.20	< 0.091	< 0.091	< 0.091	< 0.091	< 0.091		< 0.091
Dieldrin	< 0.044	< 0.044	< 0.044	<0.10	<0.10	< 0.044	<0.044	< 0.044	< 0.044	< 0.044	<0.091 <0.044	< 0.041
Endosulfan 1	1	10.04	10.044	<0.10	< 0.05	X0.044	*.0.044	₹0.044	V0.044	< 0.044	< 0.044	<0.044
Endosulfan II		<del> </del>		<0.10	< 0.20							
Endosulfan Sulfate	·	<b></b>		< 0.10	<0.10							
Endrin	< 0.039	< 0.039	< 0.039	<0.10	<0.10	< 0.039	40.000	40.070	40.005			
Endrin Aldehyd	10.037		<b>~0.039</b>	<0.10	<0.10	6.039	< 0.039	< 0.039	< 0.039	< 0.039	< 0.039	< 0.039
Heptachlor	<0.040	< 0.040	< 0.040	< 0.10	<b>\$</b> <0.10	< 0.040						
Heptachlor Epoxide	<0.032	< 0.032	<0.032				< 0.040	< 0.040	< 0.040	< 0.040	<0.040	< 0.040
Methoxychlor	<0.176	< 0.176	< 0.176	<b>∌&lt;</b> 0.10	< 0.10	< 0.032	< 0.032	< 0.032	< 0.032	< 0.032	< 0.032	< 0.032
Toxaphene	<0.086	<0.026	<0.086	<0.10	< 1.0	<0.170	< 0.175	< 0.176	< 0.176	< 0.176	< 0.176	< 0.176
Aroclor-1016	< 0.050	< 0.050	< 0.050	<2.0	<2.0	< 0.086	< 0.086	< 0.086	₹0.086	< 0.086	< 0.085	< 0.086
Aroclor-1221	< 0.054	<0.050	< 0.054			< 0.050	< 0.050	< 0.050	< 0.050	< 0.050	< 0.050	< 0.050
Arocior-1232	< 0.065	< 0.065	< 0.054			< 0.054	< 0.054	< 0.054	< 0.054	< 0.054	< 0.054	< 0.054
Aroclor-1242	< 0.065	< 0.065				< 0.065	< 0.065	< 0.065	<0.765	< 0.065	< 0.065	< 0.065
Aroclor-1248	<0.090		< 0.065			< 0.065	< 0.065	< 0.065	< 0.065	< 0.065	< 0.065	< 0.063
Aroclor-1254	<0.100	< 0.090	< 0.090			< 0.690	< 0.030	< 0.090	< 0.090	< 0.090	< 0.020	< 0.090
Aroclor-1260	< 0.100	< 0.100	<0.100			< 0.100	<.0.100	< 0.100	< 0.100	< 0.100	< 0.100	< 0.100
Aroclor-1268		< 0.100	< 0.100			< 0.100	< 0.100	< 0.100	< 0.100	< 0.100	< 0.100	< 0.100
PCB-1016/1242	< 0.100	< 0.100	< 0.100			< 0.100	< 0.100	< 0.100	< 0.100	< 0.100	< 0.100	< 0.100
PCB-1221	<del> </del>			<1.0	<1.0							
PCB-1232	<del> </del>			<1.0	<1.0							
PCB-1248	<del> </del>			<1.0	<1.0							
PCB-1254	<del></del>			< 1.0	<1.0							
PCB-1250				<1.0	<1.0							
PCB-1268	<del> </del>			<1.0	< 1.0							
2,4-D	<0.29	< 0.29										
2,4-D Silvex(2,4,5-TP)	< 0.29		< 0.29	<2.0	<2.0	< 0.29	< 0.29	< 0.29	< 0.29	< 0.29	< 0.29	< 0.29
		< 0.34	< 0.34	< 2.0	<2.0	< 0.34	<0.34	< 0.34	< 0.34	< 0.34	< 0.34	< 0.34
Aldicarb	< 0.50	< .50	< .50	< 10	< 10	<0.50	< .50	<.50	< 0.50	< .50	<.50	<0.50
Carbofuran	<5.0	< 5.0	< 5.0	< 10	< 10	< 5.0	< 5.0	< 5.0	< 5.0	< 5.0	< 5.0	< 5.0
1,2-Dibromo-3-Chloropropane	<0.10	< 0.10	< 0.10			< 0.10	< 0.10	< 0.16	< 0.10	< 9.10	< 0.10	< 0.10
Ethylene dibromide	< 0.10	< 0.10	< 0.10			< 0.10	<0.10	< 0.10	< 0.10	< 0.10	< 0.10	< 0.10
Trifluralin	II	i					<u> </u>					1

Parameter	SPRING	SPRING
E DI MIRICE	9/9/99	11/1/99
Acenaphthene		
Acenaphthylene		
Alachlor (lasso)	<1.0	< 1.0
Atrazine	<1.0	<1.0
Anthracene		
Eenzidine		
Benzo(a) anthracene Benzo(b) fluoranthene		
Benzo(k)fluoranshene		
Berzo(a)pyrene		
Benzo(g,h,i)perylene		
Benzyl alcohol	-,	
Benzyl butyl phthalate		
Bis(2-chloroethyl) ether		
Bis(2-chlorethoxy)methane		
Bis(2-cthylhexyl)phthalate	<1.0	<1.0
Bis(chloromethyl)ether	<1.0	<1.0
Bis(2-chloroisopropyl)ether	<del> </del>	
4-Bromophenyl phenyl ether 4-Chloroaniline		
2-Chloronaphthalene		
4-Chlorophenylphenyl ether		
Chrysene		
Dibenzo(a,h) anthracene		
Di-n-butylphthalate	< 1.0	<1.0
1,2-Dichlorobenzene		
1,3-Dichlorobenzene		
1,4-Dichlorobenzene		
3,3-Dichlorobenzidine	ļ	
Diethyl phthalate	<1.0	<1.0
Dimethyl phthalare	<1.0	< 1.0
2,4-Dinitrotoluene 2,6-Dinitrotoluene		
Di-n-octylphthalate		
Fluoranthene		
Fluorene		
Hexachloro-1,3-butadiene		
Hexachlorocyclopentadiene		
Hexachloroethane		
Indeno(1,2,3-cd)pyrene	ļ	
Isophorone	<1.0	< 1.0
2-Methylnaphthalene		
Naphthalene	<1.0	<1.0
2-Nitroaniline 3-Nitroaniline	<del></del>	
4-Nitroaniline	<del> </del>	
Nitrobenzene		
N-Nitrosodimethylamine	<del> </del>	
N-Nitrosodiphenylamine		
N-Nitrosodi-n-propylamine		
Phenanthrene		
Pyrene		
Pyridine	ļ	ļ
1,2,4-Trichlorobenzene		
Benzolc Acid	ļ <u>.</u>	<b> </b>
4-Chloro-3-methylphenol		ļ
2-Chlorophenol	<u> </u>	ļ
p-Cresol	<1.0	<1.0
Cresols, Total		<del> </del>
2,4-Dichlorophenol		<del></del>
2,4-Dimethylphenol 2,4-Dinitrophenol		<del>-</del>
er - magnitud		L

Parameter	SPRING 9/9/99	SPRING 11/1/99		
2-Methyl-4,6-dinitrophenol				
2-Nitrophenol				
4-Nitrophenol				
Parathion	<1.0	<1.0		
Pentachlorophenol	<1.0			
Phenol	<1,0	<1.0		
2,4,5-Trichlorophenol				
2,4,6-Trichlorophenol				
Aldrin	< 0.034	< 0.034		
alpha-BHC				
beta-BHC				
gamma-BHC	< 0.025	< 0.025		
delu-BHC				
Chlordane	< 0.037	< 0.037		
4,4'-DDD	10.051	10,037		
4,4'-DDE				
4,4'-DDT	< 0.091	< 0.091		
Dieldrin	< 0.044			
Endosulfan I	1	- 30.044		
Endosulfan II				
Endosulfan Sulfate				
Endrin	< 0.039	< 0.039		
	<b>CU.039</b>	<u> </u>		
Endrin Aldehyde	50.010	-0.042		
Heptachlor	< 0.040			
Heptachlor Epoxide	< 0.032			
Methoxychlor	< 0.176			
Toxaphene	<0.086	-		
Aroclor-1016	< 0.050			
Aroclor-1221	< 0.054			
Aroclor-1232	< 0.065			
Aroclor-1242 Aroclor-1248	<0.065	1		
	<0.090			
Aroclor-1254	<0.100			
Aroclor-1260	<0.100			
Aroclor-1268	< 0.100	< 0.100		
PCB-1016/1242	<del> </del>	<del> </del> -		
PCB-1221 PCB-1232	<del> </del>	<del> </del>		
PCB-1248	<del> </del>	<del> </del>		
FCB-1254	<del> </del>			
PCB-1260	<b> </b>	<del> </del>		
PCB-1268	<del> </del>			
2,4-D	Z0 20			
	<0.29			
Silvex(2,4,5-TP)	< 0.34			
Aldicarb	<.50			
Carbofuran	< 5.0			
1,2-Dibromo-3-Chloropropane Ethylene dibromide	<0.10	7		
	< 0.10	< 0.10		
Trifluralin	l	L		

Attachment 5

Boring Logs

**Dixon Marquette Cement Company** 

Monitoring Well MW1-D (east side of CKD landfill)

Drilled 19-March-98 by Geotechnical Services, Inc., Davenport, Iowa

Continuous rock core sampling (2 inch diameter by 5 feet long) with 4.25 inch air hammer Total depth 78 feet

Depth of Top	Thickness	Unit	Description
		Pecatonica Formation	
		Medusa member	
0	4		Limestone, yellow with gray dolomitic mottling, rare brachiopod
		New Glarus member	
4	9		Limestone, yellow with dark yellow mottling, some brachiopods, gastropods
		Dane member	
15	4		Limestone, yellow with large dolomitic or shaly mottling, some brachiopods, pyritic at 19 feet
19	8.5		Dolomite, gray, sandy, vuggy from 19 to 29 feet
		Chana member	
27.5	8		Dolomite, gray, sandy, pyritic
		Glenwood Formation	
35.5	3		Sandstone, fine grained quartz, pyritic
38.5	4		Siltstone, gray, fissile
42.5	1.5		Sandstone, quartz, some pyrite
		St. Peter Formation	
44			Sandstone, quartz, poorly c emented

## Dixon Marquette Cement Company

## Monitoring Well MW2-D (south side of CKD landfill)

Drilled 16-March-98 by Geotechnical Services, Inc., Davenport, Iowa

Continuous rock core sampling (2 inch diameter by 5 feet long) with 4.25 inch air hammer Total depth 60 feet

Depth of Top	Thickness	Unit	Description					
0	1.5		Fill material, crushed rock					
		Pecatonica Formation						
		Medusa member						
1.5	4.5		Limestone, yellow with gray dolomitic mottling, rare brachiopod					
		New Glarus member						
6	1		Limestone, yellow with dark yellow mottling					
7	4.5		Limestone, yellow with gray dolomitic mottling, some brachiopods, gastropods, crinoids					
11.5	4.5		Limestone, yellow and gray, stylolites at 12 feet					
		Dane member						
16	7.5		Limestone, yellow with large dolomitic or shaly mottling, some brachiopods, pyritic at 19.5 feet, calcite in vig. 122 feet					
23.5	<b>4</b>		Dolomite, gray, sandy, vuggy from 25 to 26 feet and 27 to 27.5 feet, calcite in vug at 24 feet, some pyrite					
27.5	1.5		Limestone, yellow and gray with dolomitic mottling					
		Chana member						
29	7.5		Dolomite, gray, sandy, pyritic, vuggy at 32 to 35.5 feet					
		Glenwood Formation						
36.5	2.5		Sandstone, fine grained quartz, some pyrite					
39	0.5		Siltstone, sandy, gray, some pyrite					
39.5	1.5		Siltstone, gray, fissile					
41	2		Sandstone, quartz, some pyrite					
43	1		Siltstone, gray, fissile					
44	3		Sandstone, quartz, mottled, some pycite					
		Et. Peter Formation	•					
47	2.5		Sandstone, quariz, poorly cemented, pyritic at 47 feet					
49.5	10		Sandstone, quartz					

**Dixon Marquette Cement Company** 

Monitoring Well MW3-D (west side of CKD landfill)

Drilled 17-March-98 by Geotechnical Services, Inc., Davenport, Iowa

Continuous rock core sampling (2 inch diameter by 5 feet long) with 4.25 inch air hammer Total depth 78 feet

Depth of Top	Thickness	Unit	Description
0	1.5		Fill material, crushed rock
		Mifflin Formation	
1.5	8.5		Limestone, yellow, fossiliferous, brachiopods
10	8.5		Limestone, yellow, mottled
		Pecatonica Formation	
		Medusa member	
18.5	4.5		Limestone, yellow with gray dolomitic mottling, rare brachiopod
		New Glarus member	•
23	11.5		Limestone, yellow with dark yellow mottling, some brachiopods, and gastropods
		Dane member	
34.5	13		Limestone, yellow with large dolomitic or shaly mottling, some brachiopods, pyritic at 19.5 feet, calcite in vug at 22 feet
		Chana member	
47.5	7.5		Dolomite, gray, sandy, pyritic
		Glenwood Formation	
55	2.5		Sandstone, fine grained quartz, some pyrite
57.5	1.5	•	Siltstone, gray, fissile
59	2		Sandstone, quartz, some pyrite
61	2		Siltstone, gray, fissile
63	3		Sandstone, quartz, mottled, some pyrite
		St. Peter Formation	
66			Sandstone, quartz, poorly cemented

#### **Dixon Marquette Cement Company**

#### Monitoring Well MW4-D (northwest side of CKD landfill)

Drilled 18-March-98 by Geotechnical Services, Inc., Davenport, Iowa

Continuous rock core sampling (2 inch diameter by 5 feet long) with 4.25 inch air hammer Total depth 60.5 feet

Depth of Top	Thickness	<u>Unit</u>	Description					
0	8.5		Fill material, crushed rock					
		Pecatonica Formation						
8.5	20.5		Limestone, yellow and gray with dolomitic mottling					
		Chana member						
29	7.5		Doiomite, gray, sandy, pyritic, vuggy					
		Glenwood Formation						
38.5	2.5		Sandstone, fine grained quartz, some pyrite					
41.5	4.5	Siltstone, gray, fissile						
46	2		Sandstone, quartz, mottled, some pyrite					
		St. Peter Formation						
48			Sandstone, quartz, poorly cemented					

#### Boring Log Dixon Marquette Cement Company Leachate Well LW-3

Drilled 07-July-99 by Geotechnical Services, Inc., Davenport, Iowa Hollow stem auger to 18 feet deep. 4.25 inch air hammer to 42 feet deep. Total depth 42 feet

Depth of Top	Thickness	<u>Unit</u>	Description
0	15	·	Brown moist rocky soil
		Mifflin?	
15	3		Rock, difficult drilling
	F	Pecatonica Formation	
32	10		Yellow brown dolomite, wet

**Dixon Marquette Cement Company** 

#### Monitoring Well MW5-D (west of CKD landfill along White Oak Lane)

Drilled 1-June-99 by Geotechnical Services, Inc., Davenport, Iowa

Continuous rock core sampling (2 inch diameter by 5 feet long) with 4.25 inch air hammer Total depth 49.5 feet

Depth of Top	Thickness	Unit	Description					
0	1		Top soil, brown, organic material					
1	1 -		Fill material					
2	1		Brown clay, trace silt, dry					
3	3		Red (brick color) clay, crumbles easily					
6	2		Light brown silt w/ trace sand, moist, gray mottles, re l/rust stains					
8	2		Brown to gray sandy silt some limestone chips					
10	12		Brown sand trace gravel					
22	4		Well graded brown soil, poorly sorted					
26	7		Brown sand with trace gravel					
33	1	Co Barrier Consider	Gray, clayey sandstone					
34	5	St. Peter Formation	Gray sandstone with brown mottling					
39	5		Gray to yellow/white sandstone					
44	5.5		Yellow to gray/green sandstone					

**Dixon Marquette Cement Company** 

Monitoring Well MW6-D (west of CKD landfill along White Oak Lane)

Drilled 7-June-99 by Geotechnical Services, Inc., Davenport, Iowa

Continuous rock core sampling (2 inch diameter by 5 feet long) with 4.25 inch air hammer Total depth 40 feet

Depth of Top	Thickness	<u>Unit</u>	Description
0	1		Top soil, brown, organic material
1	1		Brown clay with organic material
2	2.75		Brown clay with trace sand with rust and yellow mottles
4.75	7.75		Brown sand with limestone chips, and trace silts
12.5	0.5		Gray sediment/silty material
13	9		Brown sand with trace gravel, brown, fine. trace silt
22	0.5		River Rock
22.5	4		Brown sand with gravel and trace silt
26.5	0.5		Light brown clay
27.5	6.5	St. Peter Formation	Light gray to gray sandstone with sand lens, stains, and brown mottling
34	. 6		Gray to green sandstone with brown mottling

**Dixon Marquette Cement Company** 

Monitoring Well MW7-D (west of CKD landfill along White Oak Lane)

Drilled 9-June-99 by Geotechnical Services, Inc., Davenport, Iowa

Continuous rock core sampling (2 inch diameter by 5 feet long) with 4.25 inch air hammer Total depth 34.5 feet

Depth of Top	Thickness	Unit	<u>Description</u>
0	0.5	•	Top soil, brown, organic material
		Pecatonica Formation	
0.5	1.5		Gray to light brown limestone with brown mottles
3	6		Brown to gray limestone with fossils
9	4		Gray to dark gray limestone signs of dolomite
• 13	2		Light brown limestone with rust stains
		Glenwood Formation	
15	6		Dark gray shale with greasy feel Fracture noted from 15.5' to 16.5'
		St. Peter Formation	
21	13.5		Green to gray sandstone

**Dixon Marquette Cement Company** 

Monitoring Well MW8-D (west of CKD landfill along White Oak Lane)

Drilled 11-June-99 by Geotechnical Services, Inc., Davenport, Iowa

Continuous rock core sampling (2 inch diameter by 5 feet long) with 4.25 inch air hammer Total depth 39.5 feet

Depth of Top	Thickness	Unit	Description
0	. 1		Top soil, brown, organic material with white limestone chips
		Pecatonica Formation	
1	4		White limestone with brown mottles and fossilized material
5	6		Brown limestone with gray to dark gray mottles
11	1		Quartz or dolomite material with silica traces
12	1		Fossils observed in stone
13	2		Limestone, gray exterior with brown interior
15	3		White limestone with reddish brown mottles
18	2		Transition zone: alternating layers of Glenwood shale and Pecatonica limestone.
20-	1		Clay with cust mottles
21	0.5		Blue/gray greasy clay
21.5	3		Dark gray shale
24.5	2.5		No recovery from boring
		St. Peter Formation	
27	0		Sandstone observed from cuttings
34	3		Friable sandstone
37	2.5		Solid sandstone

	Illinois Environmental Protect	io	n	Age	ncy		Fi	eld	В	orii	ng Log room   of
Sits ille H	le.:CountyLel				Borin	g No	. W	WI	-5		Nonitor Well No. MWI-S
Site File N	Dixon Marquette Cement Co.										
Fed. ID. N											Rotory Depth
Quadrangle	Dixon East Sec 27 1. 22N	R.	91	Ε	Date:	Stor	t				Finlan 3-20-98
UTM (or S Plane) Cod	Ricte prd. H.(X) E(Y)										•
Latitude	Longitude						AM.	PLE			Personnel
	tion Eside of landfill					끭	8	25.	N Si	2	G - Mark Zell O - Mott White
Orliling Equ	ulpment 41 air haromer	phic	69	Depth In feet	SAUPLE NO.	SAMPLE TYPE	SAMPLE RECOVERY (X)	PEHETROMETER	ALUES V¥ COU	OINGS	G - Mark Zell O - Matt White H - Gary Kimber H - REMARKS
Elev.	DESCRIPTION OF MATERIALS	ဦ	3	ខ្លួន	3	इ	35	ģ	ZE	\$5	REMARKS
<u> -</u>	Limestone, yellow with gray dolomitic mottling,			E							·
F	Limestone, yellow with dark yellow mottling, some brachiopods, gastropods			-							Continuous core
726.70	some oracmopous, gastropous			-10	ļ	<b> </b>	·	! 		<b></b>	County of Colf
<b>†</b>	Limestone, yellow with large dolomitic or shaly mottling, some brachiopods, pyritic at 19 feet			_							
71674				- 20		<u> </u>					
F				_							
	Dolomite, gray, sandy, pyritic										
706:70				30		ļ					
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Site File N	Dixon Marquette Cement Co.				Surfa	<b>=</b> 0	evetic	a 7	36.8	14	Completion Copth 78-ft
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Boring Loc	etion Eside of landfill				3	TYPE	8	TER	NIS)	2	G - Mark Zell O - Matt White
Orllling Equ	ulpment 44 air hummer	21467	2007	Depth In feet	SAUPLE NO.	SAUPLE TY	SAMPLE RECOVERY	PEHETROWETER	VALUES O¥ COU	NE STAN	G - Mark Zell O - Matt White H - Gary Kimler H - REMARKS
Elev.	DESCRIPTION OF MATERIALS	18	٤٤	20	3	क्र	35	Gd	zΞ	8€	REMARKS
-	Limestone, yellow with gray dolornitic mortling, rare brachiopod			<u> </u> -							
-	Limestone, yellow with dark yellow mortling, some brachiopods, gastropods			<u> </u>							Continuous core
726.84				-10 				••••		••••	
[	Limestone, yellow with large dolomitic or shaly mottling, some brachiopods, pyritic at 19 feet										
716.84	Dolomite, gray, sandy, vuggy from 19 to 29 feet			-20	ļ			• • • • • • • • •			
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706.84	Dolomite, gray, sandy, pyritic			-30		ļ					
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-696.84	Sandstone, fine grained quartz, pyritic Siltstone, gray, fissile		<b>Z</b>	_							
-670:34	Sandstone, quartz, some pyrite			—40 -		•••••	••••	•••			•
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Revised 28-May-99

Ellinois Environmental Protection Agency Field Boring Log - 1 a 1														
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Site File Name Dixon Marquette Cement Co.														
Fad. 10. N						oth								
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Ortilling Equ	ulpment 44 air hammer	phíc	ر ن	SAMPLE NO.	ונ	SAMPLE RECOVERY	PEMETROMETER	UES COU	or HNU NGS	H - Gary Kimler				
Eley.	DESCRIPTION OF MATERIALS	g 2	Depth In feet	SAMP	SAMPLE	SECON	PENE	N VAL	188 188 188	REMARKS				
	Fill material, crushed rock.  Limestone, yellow with gray dolomitic monthing	T	<del> </del>	-						CANERIAN				
	Limestone, yellow with dark yellow mottling Limestone, yellow with gray dolomitic mottling, some		F											
-690.67	brachiopods, gastropods, crinoids		F 10			<u> </u>				Continuous core				
	Limestone, yellow and gray, stylolites at 12 feet		F			.								
670.67	Limestone, yellow with large dolomitic or shaly mottling, some brachlopeds, pyritic at 19.5 feet		F											
	Dolomite, gray, sandy, vuggy, some pyrite		<b>-20</b>							9				
-	Limestone, yellow and gray with dolomitic mottling		Ē					ļ						
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650.67	Sandstone, fine grained quartz, some pyrite Siltstone, sandy, gray, some pyrite	2	-40											
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Illinois Environmental Protect	tion	Age	ncy	<del></del>	Fi	eld	В	orir	ag Log Foot
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Site File Hame Dixon Marquette Cement Co.									
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Quedrengle Discon East Sec. 27 T. 22N	R 91								
VTM (or State Plane) Coord, M.(X) E(Y)							* * *****		•
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Boring Location S of landfill			j j	TYPE	E	200	NIS)	2	0 - Mark Zell 0 - Matt White
Orilling Equipment 44 air hammer	Graphic Log	oth feet	SAUPLE NO.	SALIPLE T	אויועגא וונכוסיווי (א)	РСИБНОМЕТЕН	27U.¥	ON LIN	G - Mark Zell G - Matt White H - Gary Kinler H - REMARKS
Elev. DESCRIPTION OF MATERIALS	ادق	25	3	75	33	۲	z	85	REMARKS
Fill material, crushed rock. Limestone, yellow with gray dolomitic mottling Limestone, yellow with dark yellow mottling Limestone, yellow with gray dolomitic mottling, some brachiopods, gastropods, crinoids Limestone, yellow and gray, stylolites at 12 feet		- - - 10		-				,	Continuous core
Limestone, yellow, with large dolomitic or shaly mottling, some brachiopods, pyrice at 19.5 feet  Dolomite, gray, sandy, vuggy, some pyrite  Limestone, yellow and gray with dolomitic mottling		- 20 -9		a.	5 <sup>2</sup>				•
Dolomite, gray, sandy, pyritic, vuggy at 32 to 35.5 feet   Sandstone, fine grained quartz, some pyrite.		- 30					<u> </u>		-
Siltstone, sandy, gray, some pyrite Sandstone, quartz, some pyrite Siltstone, gray, fissile Sandstone, quartz, mottled, some pyrite! Sandstone, quartz, poorly cemented, pyritic at 47 feet Sandstone, quartz		40  - - - -50		••••					
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	Illinois Environmental Protec	tion	ı Age	ney	,	Fi	ielo	i B	ori	ng Log 🗫   «
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Site File N	isme Dixon Marquette Cement Co.					سعد العجواة	:n 6	93.9	6	Completion Cepth 54 Dt
Fed. ID. N	lo									Rotory Depth
Quedrengle	Dixon East Sec. 27 T. 221	J R								Finish 3-20-98
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Boring Loc	w side of landfill				ڀ	ક્ર	TER	(3)		G - Mark Zell
Orliling Equ	ulpment 44° air hammer	phlc	Log Depth In feet	SAUPL: NO.	SALIFICE TYPE	SAUPLE RECOVERY (X)	PENETROMETER	ALUES W COU!	OINCS	G - Mark Zell O - Matt White H - Gary Kimker H - REMARKS
Elev.	DESCRIPTION OF MATERIALS	ဦ	2 2 2	3	क्र	35	PEN	25	8₹	REMARKS
<u> </u>	Fill material, crushed rock Limestone, yellow, fossiliferous, brachiopods		E							
-			E							Continuous core
683.96	Limestone, yellow, mottled		-10				•••••			
<b>F</b>			F							
673.96	Limestone, yellow with gray dolomitic mortling, rare brachiopod		-20							
<u> </u>	Linestone, yellow with dark yellow mottling, some brachiopods, and gastropods		F							
-663.96	some enteriores, and garacters		F_					ĺ		
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<u> </u>	Limestone, yellow with large dolomitic or shaly mottling, some brachiopods, pyritic at 19.5 feet,		F						ļ	
65396	calcite in vug at 22 feet		-40							
<u> </u>			F							,
643.46	Dolomite, gray, sandy, pyritic		F_							
			-50		•••••	······		·····†		
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	llinois Environmental Protect	io	n	Age	псу		Fi	eld	B	orii	ng Log 🗫 📗 🗸 📗			
Sits File No.: County Lee						Saring No. MW3-D Wanitar Wall No. MW3-								
Site File Name Dixon Marquette Cement Co. S						Surfaces Devotion 640.26 Completion Cooth 78 ft								
Fed. ID. N											Rotory Depth			
Quedrengla	Dixon East Sec 27 T. 22N	. ጼ	9	E	Octes	Stor	t				Finish 3-20-98			
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Baring Lac	ation Wside of landfill				Š.	35	8	EJE S	Z S	2	G - Mark Zell G - Matt White			
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Elev.	DESCRIPTION OF MATERIALS	٤	2 ا	20.5	SAUPLE	র	35	Ġ	ΣĒ	8€	REMARKS			
	Fill material, crushed rock Limestone, yellow, fossiliferous, brachiopods			-  -  -							Continuous core			
-680.26	Limestone, yellow, mottled			<u>-10</u>										
-670.26	Limestone, yellow with gray dolomitic mottling,			-			·							
-	rare brachiopod Limestone, yellow with dark yellow monling,			20  				••••						
-66026	some brachiopods, and gastropods			- 4:										
-				30 -						••••	:			
-650.26	Limestone, yellow with large dolomitic or shaly mottling, some brachiopods, pyritic at 19.5 feet, calcite in vug at 22 feet			-										
-				40 				•••••						
<u> </u>	Dolomite, gray, sandy, pyritic			_						İ				
640.26				S0	ļ									
-630.26	Sandstone, fine grained quartz, some pyrite Siltstone, gray, fissile	\ \ \		-										
- 22.20	Sandstone, quartz, some pyrite Siltstone, gray, fissile			40 	•••••	••••				•••••				
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Revised 28-May-99

Ellinois Environmental Protection Agency Field Boring Log - 1 0										
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Boring Location	NW of landfill			ا ا	34	શ	TER	NIS)	Э	G - Mark Zell O Matt White
Ortiling Equipm	nent 44° air hammer	phlc	Log Depth In feet	SULPLE NO.	SALIPLE TYPE	SAUPLE RECOVERY (X)	PENETROWETER	WLUES DW COU	DINCS	G - Mark Zell O Matt White H Gary Kimler H  REMARKS
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	Il material, crushed rock mestone, yellow and gray with dolomitic mottling		7-1-1							Confinuous core
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672.44			-20						••••••	
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GCATAA DO	plomite, gray, sandy, pyritic, vuggy		-30							
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	ndstone, fine grained quartz, some pyrite	Ц								
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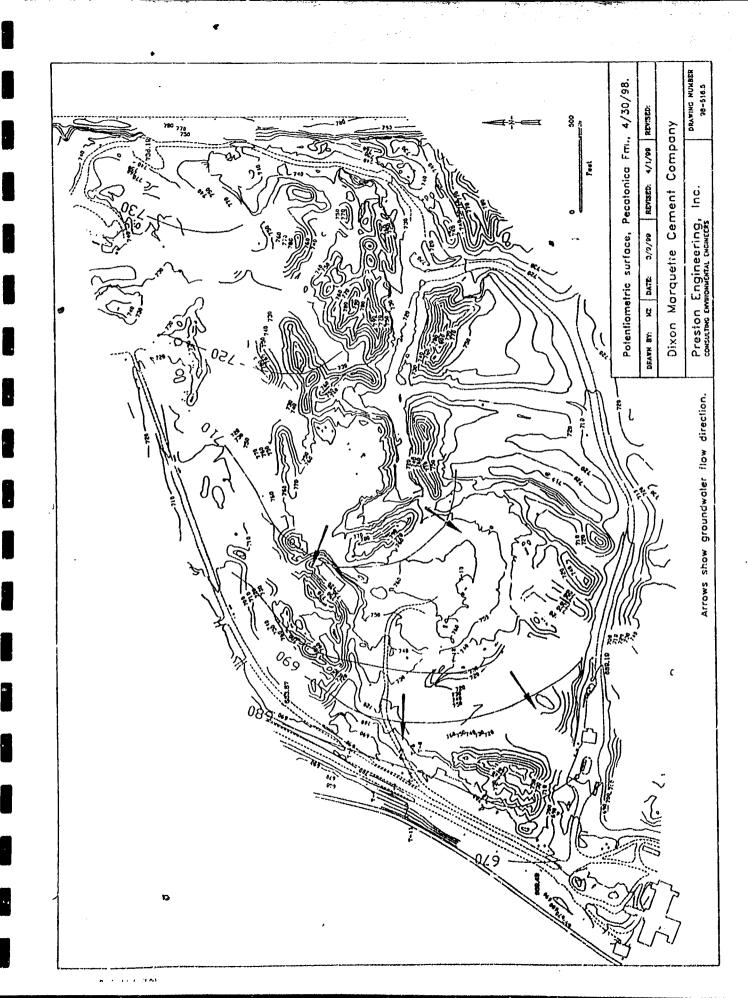
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Sits File No	c.: County Lee		····	∄c <i>r</i> i	ng No	<u>M</u> 1	W4.	- D		Monitor Well No. MW4-D
Fed. ID. N	Dixon Marquette Cement Co.									
										Ratery Depth
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Plene) Coc	rd. н.(X) E(Y)			<b>,</b>	····					•
	Longituda		<u> </u>			AM	PLE	S	<b></b>	Personnel
Boring Lac	otion NW of landfill			- Ş.	ير	E	200	(SIMES)	5	G - Mark Zell G - Matt White
Orilling Equ	DESCRIPTION OF MATERIALS	Plc	uphic John Sth feet		SAUPLE TYPE	ַבָּ קַנְיַי נָ	PENETROMETUR	N VALUES	PC S	0 - Matt White н - Gary Kimler н -
Elev.	DESCRIPTION OF MATERIALS	5	Log Depth	SAMPLE SAMPLE	Sku	אוורכסענוגע (א)	PUNE	35	OXY of RESOUR	REMARKS
F	Fill material, crushed rock	$\prod$	-	1	†					AMMANA
F			F							Continuous core
682.69	Limestone, yellow and gray with dolomitic mortling		-10	,				,,,,,,		CONTINUOUS COLE
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-662.69	Dolomite, gray, sandy, pyritic, vuggy		-3		ļ					· ·
			F		Ì					
-652.41	Sandstone, fine grained quartz, some pyrite		<u>ا</u> ۔							
<u> </u>	Siltstone, gray, fissile	/	-4	)		··· ·· <del>i</del>	••••			•
<u> </u>	Sandstone, quartz, mottled, some pyrite	-	+							
642.69	Sandstone, quartz, poorly cemented	,	/F-50	·						
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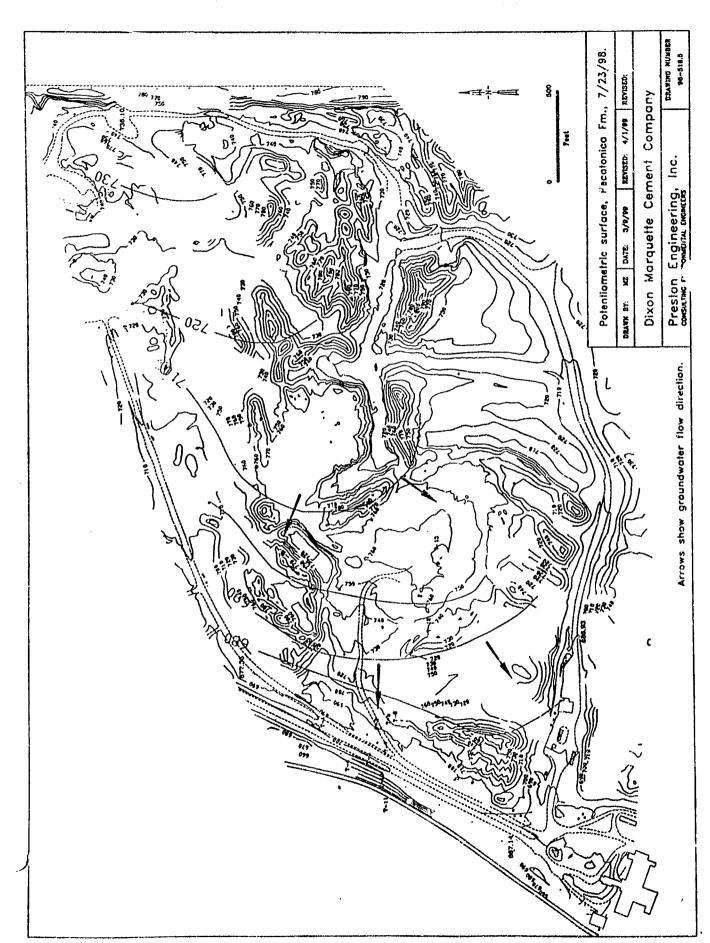
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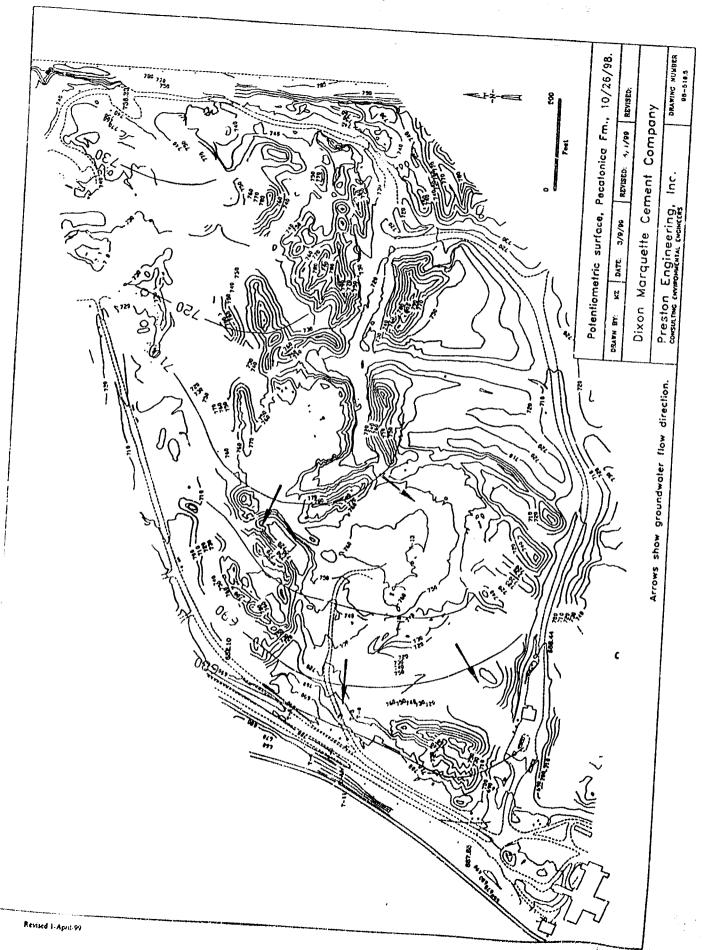
Revised 28-May-99

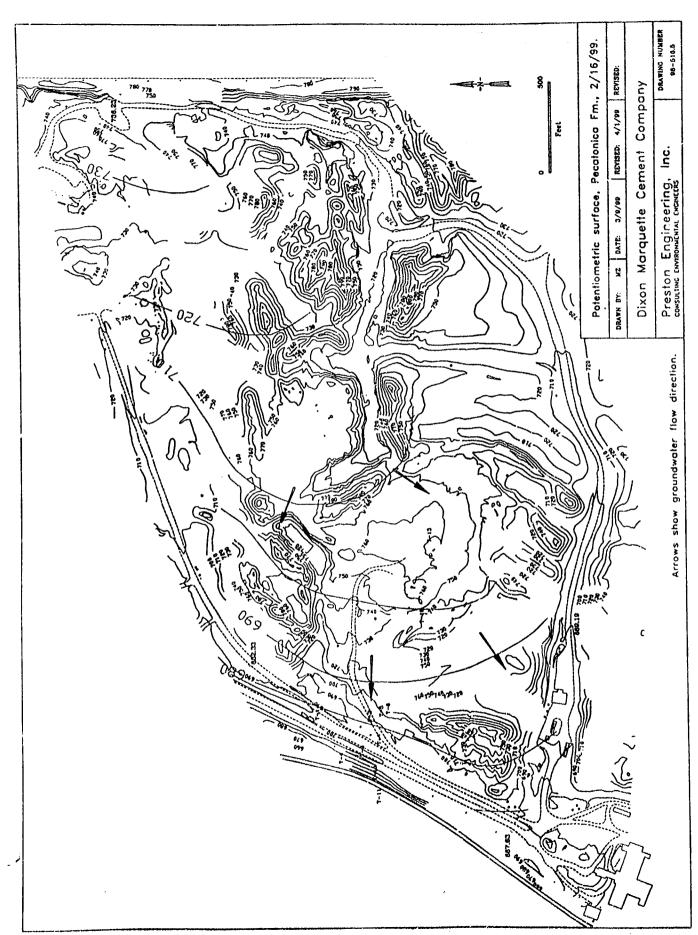
Attachment 6

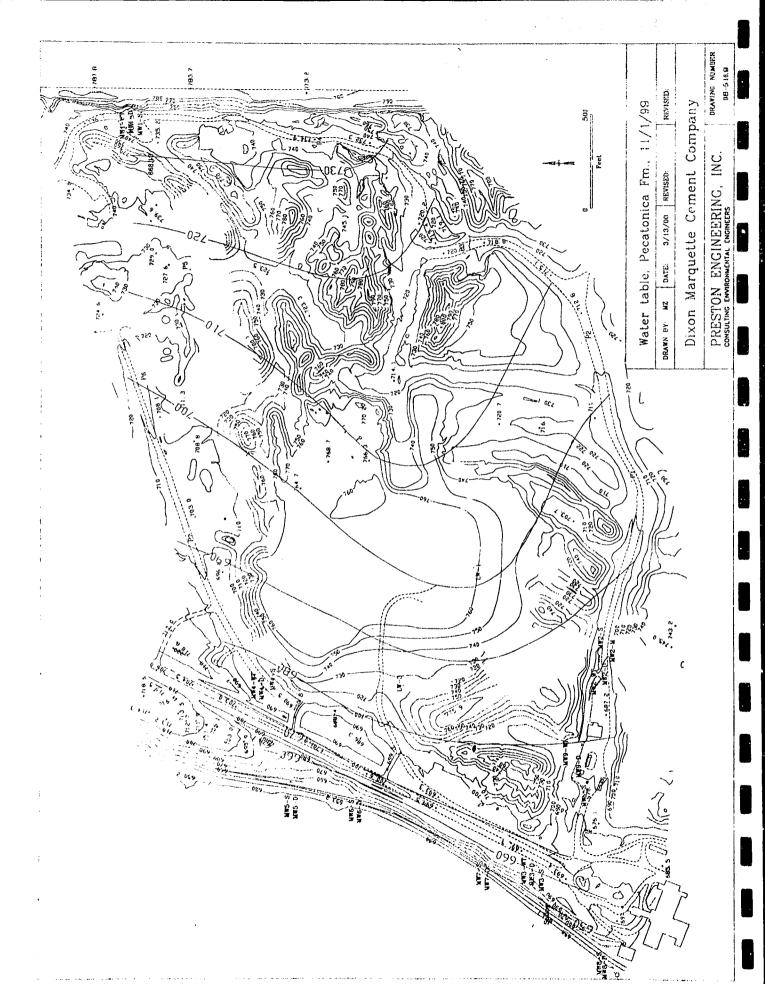
Potentiometric Surface Maps

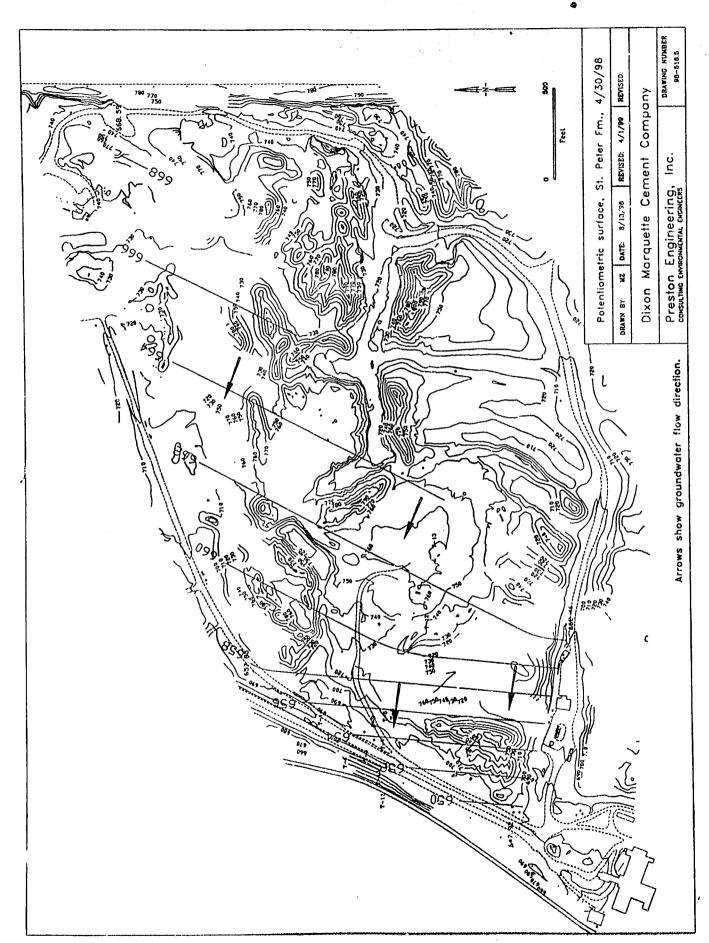


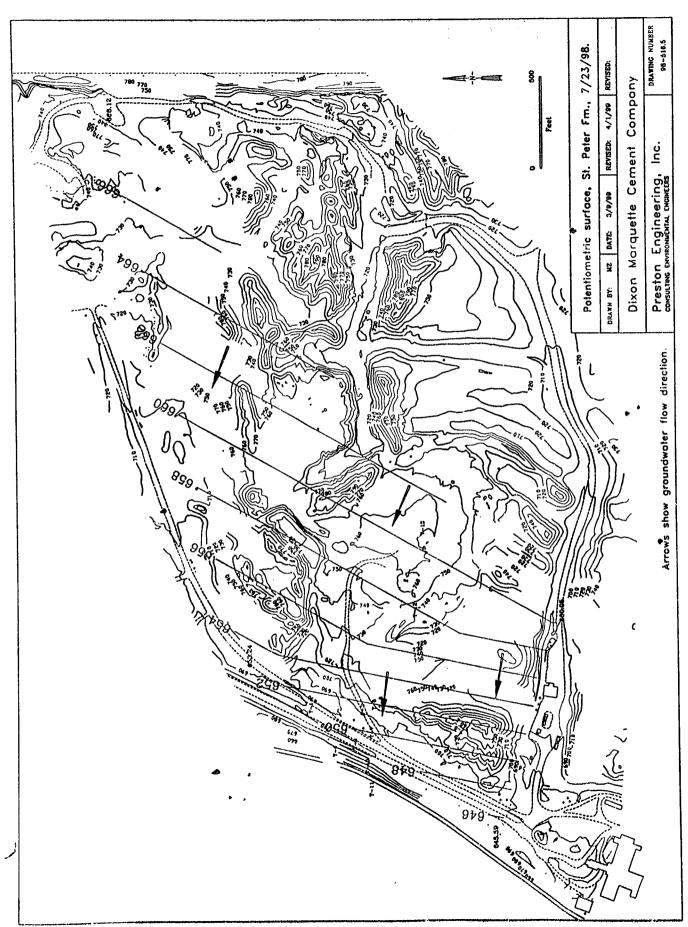


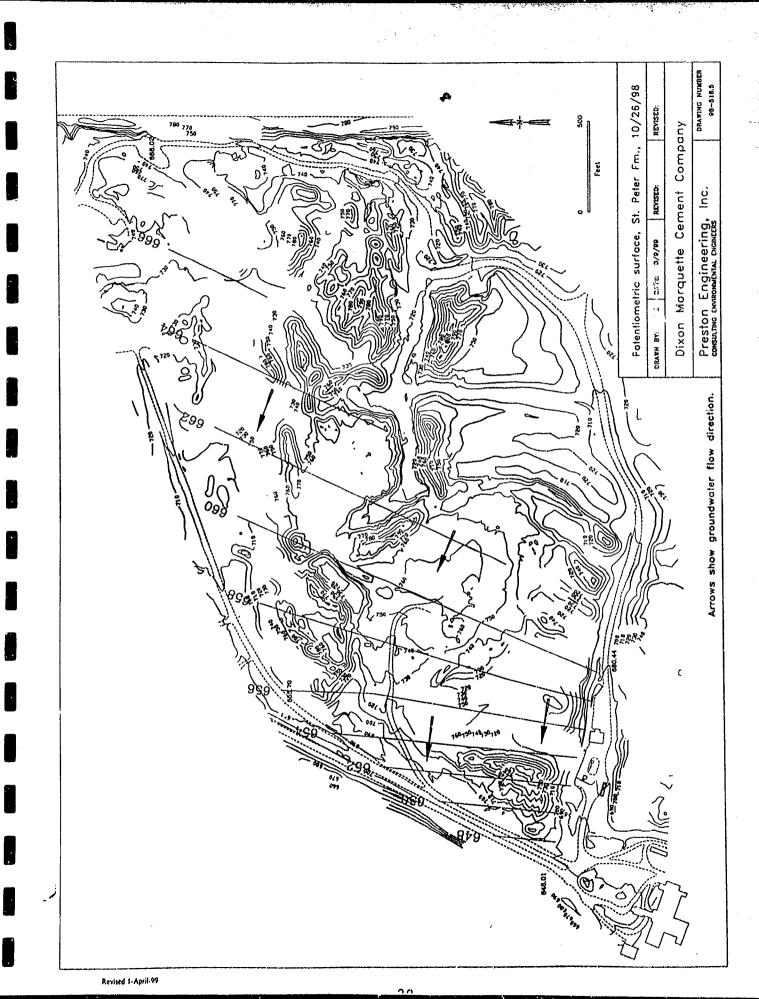




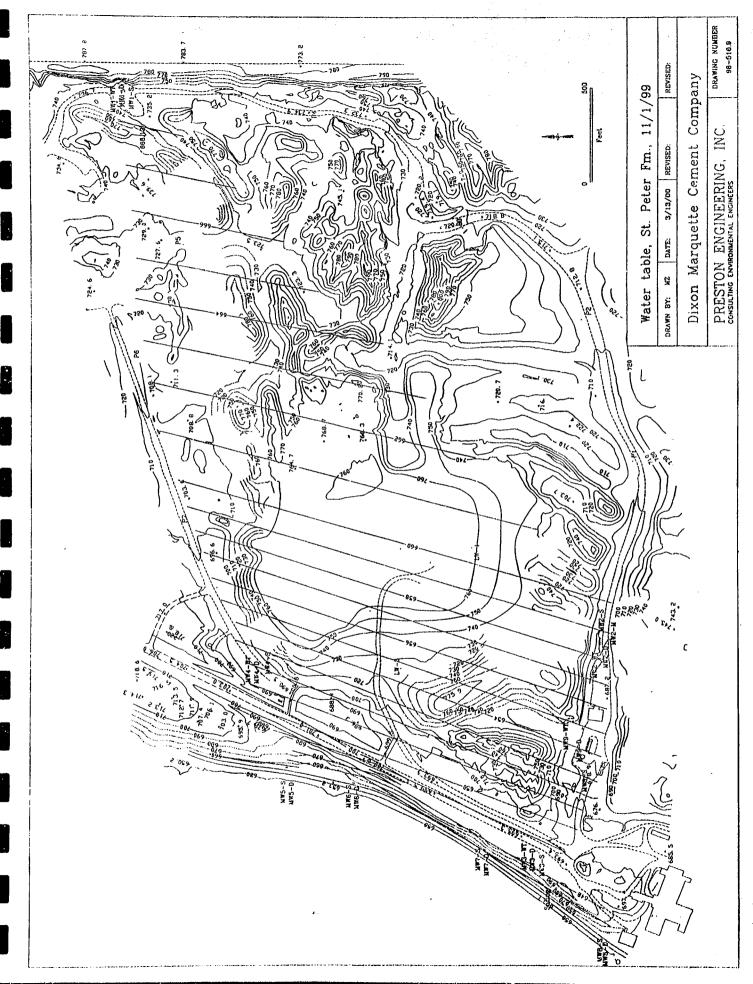






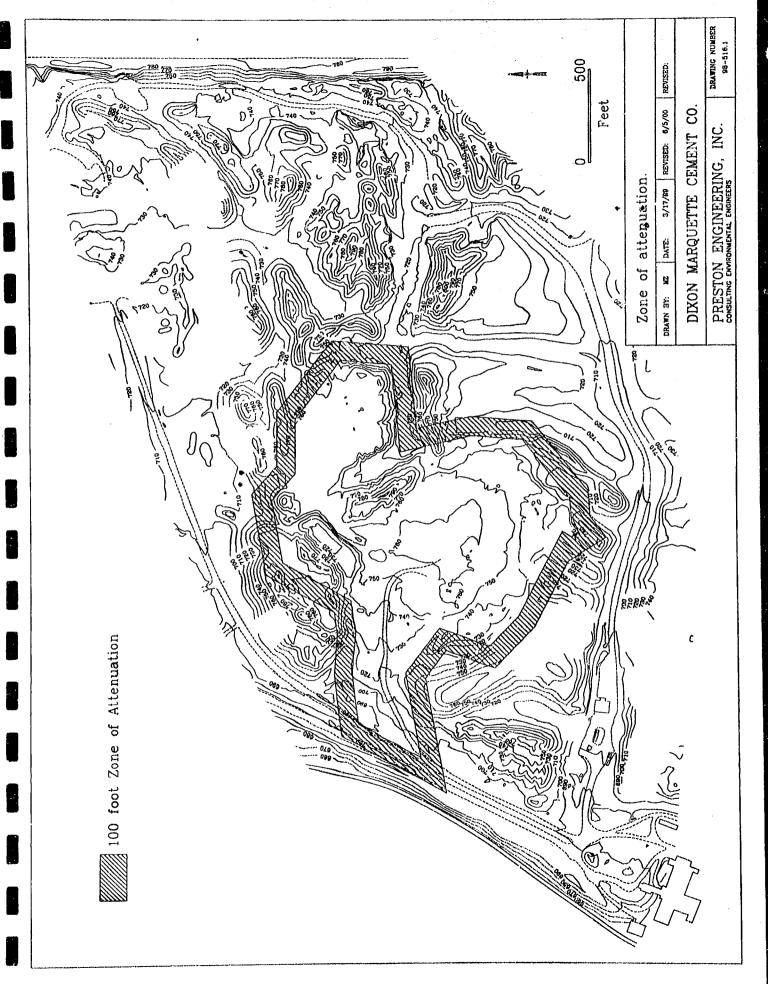






Attachment 7

Zone Of Attenuation



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Preston Engineering, Inc.

environ mental consultants

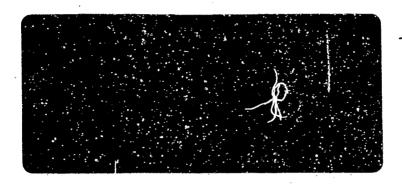
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STATE OF ILLINOIS
Pollution Control Board



A501-10

#### PETITION OF DIXON MARQUETTE CEMENT COMPANY FOR ADJUSTED STANDARD FROM: 35 ILL. ADM. CODE PARTS 811 & 814

**EXHIBITS 6-10** 

March 15, 2001

Prepared by:

Preston Engineering, Inc. 4436 North Brady-Street Davenport, Iowa 52806 (319) 388-8288- phone (319) 388-9003- fax



## ILLINOIS ENVIRONMENTAL PROTECTION AGENCY

1021 NORTH GRAND AVENUE EAST, P.O. BOX 19276, SPRINGFIELD, ILLINOIS 62794-9276

217/782-0610

THOMAS V. SKINNER, DIRECTOR

September 13, 1999

Dixon-Marquette Cement Company 1914 White Oak Lane Post Office Box 467 Dixon, Illinois 61021

Re: Dixon-Marquette Cement Company

Dixon Plant

NPDES Permit No. IL0003514

Modification of NPDES Permit (After Public Notice)

#### Gentlemen:

The Illinois Environmental Protection Agency has reviewed the request for modification of the above-referenced NPDES Permit and issued a public notice based on that request. The final decision of the Agency is to modify the Permit as follows:

Outfall 003A has been redesignated Outfall 015.

Internal Outfalls 001A, 001B, 001C, 002A, 002B, 002C, 002D, 002E, 003A, 003B, 004A and 004B have been designated.

Outfall 014 has been added to the permit.

Outfalls 007, 008 and 009 have been removed from the permit due to coverage under NPDES Permit ILG840104.

Enclosed is a copy of the modified Permit. You have the right to appeal this modification to the Illinois Pollution Control Board within a 35 day period following the modification date shown on the first page of the permit.

Should you have any question or comments regarding the above, please contact David Ginder of my staff.

Very truly yours,

Thomas G. McSwiggin, P.E.

Manager, Permit Section

Division of Water Pollution Control

TGM:BJY:DPG:98030407.dlk

Attachment: Modified Permit

cc: Records

Compliance Assurance Section; Willet, Hoffman & Associates

#### Illinois Environmental Protection Agency

**Division of Water Pollution Control** 

1021 North Grand Avenue East

Post Office Box 19276

Springfield, Illinois 62794-9276

## NATIONAL POLLUTANT DISCHARGE ELIMINATION SYSTEM

Modified (NPDES) Permit

Expiration Date: November 30, 2000

Issue Date: December 4, 1995

Modification Date: September 13, 1999

Name and Address of Permittee:

Facility Name and Address:

Dixon-Marquette Cement Company 1914 White Oak Lane

Post Office Box 467 Dixon, Illinois 61021 Dixon Plant 1914 White Oak Lane Dixon, Illinois 61021

(Lee County)

Discharge Number and Name:

Receiving Waters

001A Internal Outfall

001B Internal Outfall

001C Internal Outfall

001 Combined flows from Outfalls 001A - 001C, and Storm Water Rock River

002A Internal Outfall

002B Internal Outfall

002C Internal Outfall

002D Internal Outfall

002E Internal Outfall

002 Combined flows from Outfalls 002A - 002E, and Storm Water

003A Internal Outfall 003B Internal Outfall

003 Combined flows from 003A - 003B, and Storm Water

004A Internal Outfall

004B Internal Outfall

004 Combined flows from Outfalls 004A - 004B, and Storm Water

runoff from an Inactive Limestone Mining Area and Cement

Kiln Dust Landfill

005 Storm Water Runoff from Inactive Limestone Mining Area, Aggregate Wash Plant Sedimentation Pond Overflow and

Coal Stock Pile Runoff

006 Storm Water Runoff

Storm Water Runoff from an Inactive Limestone Mining Area

and Cement Kiln Dust Landfill

015 Sawage Treatment Plant Effluent Rock River

**Rock River** 

Unnamed Tributary to the Rock River

Rock River

Rock River

Rock River

Rock River

In compliance with the provisions of the Linois Environmental Protection Act, Title 35 of III. Adm. Code, Subtitle C and/or Subtitle D, Chapter 1, and the Clean Water Act (CWA), the above named permittee is hereby authorized to discharge at the above location to the above-named receiving stream in accordance with the standard conditions and attachments herein.

### Page 2

Permittee is not auth@rized to discharge after the above expiration date. In order to receive authorization to discharge beyond the expiration date, the permittee shall submit the proper application as required by the Illinois Environmental Protection Agency (IEPA) not later than 180 days prior to the expiration date.

Thomas G. McSwiggin, P.E. Manager, Permit Section
Division of Water Pollution Control

TGM:DPG:98030407.dlk

1/Month

1/Month

1/Month

1/Month

1/Month

Grab

Single Reading

Grab

Grab

Grab

Temperature\*

Sulfate

**Total Suspended Solids** 

**Total Dissolved Solids** 

#### NPDES Permit No. IL0003514

#### **Effluent Limitations and Monitoring**

CONCENTRATION LIMITS mg/l

LOAD LIMITS lbs/day DAF (DMF)

See Special Condition 5

**Monitor Only** 

**Monitor Only** 

PARAMETE	30 DAY R AVERAGE	DAILY MAXIMUM	30 DAY AVERAGE	DAILY MAXIMUM	SAMPLE FREQUENCY	SAMPLE TYPE
From the modific at all times as follows:	cation date of this permit un ws:	til the expiration dat	e, the effluent of the	following discharg	e(s) shall be monito	red and limited
Outfalls: 00	01 and 003				en en en en en en en en en en en en en e	
Flow					1/Month	Single Reading
рН	See Special Con-	dition 2			1/Month	Grab

15

30

<sup>\*</sup>At the time of each effluent sampling event, a grab sample of the cooling water intake for each outfall shall be collected and a temperature reading shall be taken and reported on the DMR form.

### **Effluent Limitations and Monitoring**

		TS lbs/day - (DMF)	CONCEN Limi'i	TRATION (S mg/l		·
PARAMETER	30 DAY	DAILY	30 DAY	DAILY	SAMPLE	SAMPLE
	AVERAGE	MAXIMUM	AVERAGE	MAXIMUM	FREQUENCY	TYPE

1. From the modification date of this permit until the expiration date, the effluent of the following discharge(s) shall be monitored and limited at all times as follows:

Outfall: 002

Flow				1/Month	Single Reading
рН	See Special Condition 2			1/Month	Grab
Temperature*	See Special Condition 5			1/Month	Single Reading
Total Suspended Solids	18.6**			1/Month	Grab
Iron	Monitor Only	3.5	7.0	1/Month	Grab
Total Dissolved Solids	Monitor Only	~		1/Month	Grab
Manganese	Monitor Only			1/Month	Grab
Sulfate	Monitor Only			1/Month	Grab

<sup>\*</sup>At the time of each effluent sampling event, a grab sample of the cooling water intake for each outfall shall be collected and a temperature reading shall be taken and reported on the DMR form.

<sup>\*\*</sup>See Special Condition 13

### Effluent Limitations and Monitoring

,	Load (.imi Daf	TS lbs/d <b>ay</b> (DMF)	CONCENTRATION LIMITS mg/l			
PARAMETER	30 DAY	DAILY	30 DAY	DAILY	SAMPLE	SAMPLE
	AVERAGE	MAXIMUM	AVERAGE	MAXIMUM	FREQUENCY	TYPE

1. From the modification date of this permit until the expiration date, the effluent of the following discharge(s) shall be monitored and limited at all times as follows:

Outfalls: 002C, 004A, 004B and 014

Flow				1/Month	Single Reading
рН	See Special Condition 2			1/Month	Grab
Total Suspended Solids		15	30	1/Month	Grab
Total Dissolved Solids**	Monitor Only			1/Month	Grab
Sulfate**	Monitor Only.			1/Month	Grab

<sup>\*\*</sup>Total Dissolved Solids and Sulfate Monitoring required for Outfall 014 only.

Flow

pН

**Total Suspended Solids** 

## NPDES Permit No. IL0003514

### **Effluent Limitations and Monitoring**

	Effluent Limitations and Monitoring						
	LOAD LIMI	TS lbs/day (DMF)	CONCENT LIMIT	TRATION 'S mg/l			
PARAMETER	30 DAY AVERAGE	DAILY MAXIMUM	30 DAY AVERAGE	DAILY MAXIMUM	SAMPLE FREQUENCY	SAMPLE TYPE	
From the modification data all times as follows:	ate of this permit un	til the expiration dat	e, the effluent of the	following discharg	e(s) shall be monitor	ed and limited	
Outfall: 004				en de la companya de		٠	
Flow					1/Month	Single Reading	
ρН	See Special Con	dition 2		e e e e e e e e e e e e e e e e e e e	1/Month	Grab	
*Temperature	See Special Con	dition 5			1/Month	Single Reading	
Total Suspended Solids**			15	30	1/Month	Grab	
Total Dissolved Solids	Monitor Only				1/Month	Grab	
Sulfate	Monitor Only				1/Month	Grab	
*At the time of each effluer reading shall be taken and **See Special Condition 1 Outfall: 002A and	d reported on the D 3.	grab sample of the MR form.	cooling water intake	e for each outfall sh	nall be collected and	a temperature	
Flow					1/Month	Singl <b>e</b> Readi <b>ng</b>	
pН	See Special Cor	idition 2			1/Month	Grab	
Total Suspended Solids			15	30	1/Month	Grab	
fron**			2.0	4.0	1/Month	Grab	
Manganese**			1.0	2.0	1/Month	Grab	
*Effluent sampling for flow **Iron and Manganese mo	v shall be continuol onitoring for Outfail	us if hardware allow 005 shall be Monit	vs otherwise it shall or only.	be once a month	single reading.		
Outfalls: 006							

35

70

See Special Condition 2

1/Month

1/Month

1/Month

Grab

Grab

<sup>\*</sup>Effluent sampling for flow shall be continuous if hardware allows otherwise it shall be once a month single reading.

## Effluent Limitations and Monitoring

	LOAD LIMI DAF	TS lbs/day (DMF)		TRATION IS mg/I		
PARAMETER	30 DAY	DAILY	30 DAY	DAILY	SAMPLE	SAMPLE
	AVERAGE	MAXIMUM	AVERAGE	MAXIMUM	FREQUENCY	TYPE

<sup>1.</sup> From the modification date of this permit until the expiration date, the effluent of the following discharge(s) shall be monitored and limited at all times as follows:

## Outfall: 015 - Sewage Treatment Plant Effluent

Flow				1/Month	
рН	See Special Condition 2			ı/Month	Grab
Total Suspended Solids		30	60	1/Month	Grab
BOD <sub>5</sub>		30	60	1/Month	Grab
Fecal Coliform	See Special Condition 1			•	ă.

#### Special Conditions

SPECIAL CONDITION 1. The daily maximum fecal coliform count shall not exceed 400 per 100 ml.

SPECIAL CONDITION 2. The pH shall be in the range 6.0 to 9.0. The monthly minimum and monthly maximum values shall be reported on the DMR form.

SPECIAL CONDITION 3. Within 180 days of the modification date of this permit, the permittee shall provide written information to the Illinois EPA on any biocides used in the Powerhouse Cooling Towers. Such information shall include the name of the chemical(s) used, active ingredients, dosage concentration, anticipated concentration in the discharge from Outfall 001A and Outfall 001, and aquatic toxicity data. The Illinois EPA may modify this permit during its term based on information provided under this Special Condition. Such modification shall follow Public Notice and opportunity for hearing.

<u>SPECIAL CONDITION 4</u>. Sampling at Outfalls 001, 002 and 003 shall be performed during dry weather periods (when no stormwater is present in the discharges).

<u>SPECIAL CONDITION 5</u>. Discharge of wastewater from Outfalls 001, 002, 003, and 004 must not alone or in combination with other sources cause the receiving stream to violate the following thermal limitations at the edge of the mixing zone which is defined by Section 302.211, Illinois Administration Code, Title 35, Chapter 1, Subtitle C, as amended:

- A. Maximum temperature rise above natural temperature must not exceed 5°F (2.8°C).
- B. Water temperature at representative locations in the main river shall not exceed the maximum limits in the following table during more than one (1) percent of the hours in the 12-month period ending with any month. Moreover, at no time shall the water temperature at such locations exceed the maximum limits in the following table by more than 3°F (1.7°C). (Main river temperatures are temperatures of those portions of the river essentially similar to and following the same thermal regime as the temperatures of the main flow of the river.)

	<u>Jan.</u>	<u>Feb.</u>	Mar.	<u>April</u>	May	June	July	Aug.	Sept.	Oct.	Nov.	Dec.
°F	60	60	60	90	90	90	90 .	90	90	90	90	60
°C	16	16	16	32	32	32	32	32	32	32	32	16

C. The monthly maximum value shall be reported on the DMR form.

SPECIAL CONDITION 6. For the purpose of this permit, Outfalls 001A-001C, 001, 002A-002E, 002, 003A-003B, 003, 004A-B, 004, 005, and 014 are designated as follows:

Outfall Number	Description
001A	Non-contact cooling water from the Powerhouse Cooling Towers
0018	Old Clay Storage Building compressor cooling water
001C	Raw Mills and Raw Mill compressor cooling water
001	Flows Tributary to Outfalls 001A, 001B, 001C and storm water
002A	Coal Pile Runoff
002B	Plant Air compressors cooling water
002C	Coal Mill cooling water
002D	Precipitator compressors cooling water, Preheater Fans cooling water, Analyzer water, Dracco Baghouse Fan cooling water, Homogenizing Silos compressors cooling water, #1, 2, 3 Kiln Piers Discharge End cooling water
002E	#1, 2, 3 Finish Mills cooling water, cement coolers, #1, 2, 3 Finish Mill compressor cooling water, #4 Finish Mill compressor cooling water, B Pump compressors cooling water
002	Flows tributary to Outfalls 002A - 002E, and storm water

#### **Special Conditions**

003A	Packhouse compressors "A" Pump cooling water
003B	Packhouse compressors instrument Air cooling water
003	Flows tributary to 003A and 003B, and storm water
004A	#1, 2, 3, Kiln Piers Feed End cooling water, #4 Piers Discharge End cooling water
0048	#4 Kiln compressors cooling water, Wheelabrator Baghouse Fan Motor Bearings cooling water, #4 Kiln Plers Feed End cooling water
004	Flows Tributary to Outfall 004A and 004B, and storm water runoff from an inactive Limestone Mining Area and Cement Kiln Dust Landfill
005	Storm water runoff from Inactive Limestone Mining Area, Aggregate Wash Plant Sedimentation Pond Overflow and Coal Stock Pile Runoff.
014	Storm Water Runoff from an Inactive Limestone Mining Area and Cement Kiln Dust Landfill

<u>SPECIAL CONDITION 7</u>. Samples taken in compliance with the effluent monitoring requirements shall be taken at a point representative of the discharge, but prior to entry into the receiving stream.

<u>SPECIAL CONDITION 8</u>. The permittee shall record monitoring results on Discharge Monitoring Report forms using one such form for each discharge each month. The completed Discharge Monitoring Report form shall be submitted monthly to IEPA, no later than the 15th of the following month, unless otherwise specified by the Agency, to the following address:

Illinois Environmental Protection Agency Bureau of Water Compliance Assurance Section 1021 North Grand Avenue East Post Office Box 19276 Springfield, Illinois 62794-9276

Flows shall be reported as a monthly average on the Discharge Monitoring Report form.

<u>SPECIAL CONDITION 9</u>. The permittee shall notify the Agency in writing by certified mail within thirty days of abandonment, cessation, or suspension of active mining for thirty days or more unless caused by a labor dispute. During cessation or suspension of active mining, whether caused by a labor dispute or not, the permittee shall provide whatever interim impoundment, drainage diversion, and wastewater treatment is necessary to avoid violations of the Act or Subtitle D, Chapter 1.

SPECIAL CONDITION 10. The Agency has determined that the effluent limitations in this permit constitute BAT/BAC for storm water which is treated in the existing treatment facilities for purposes of this permit issuance, and no pollution prevention plan will be required for such storm water. In addition to the chemical specific monitoring required elsewhere in this permit, the permittee shall conduct an annual inspection of the facility site to identify areas contributing to a storm water discharge associated with mining or industrial activities and determine whether any facility modifications have occurred which result in previously-treated storm water discharges no longer receiving treatment. If any such discharges are identified the permittee shall request a modification of this permit within 30 days after the inspection. Records of the annual inspection shall be retained by the permittee for the term of this permit and be made available to the Agency upon request.

SPECIAL CONDITION 11. Mining excavation operations shall maintain a minimum setback of 200 feet from the two private potable wells located in Sections 15 and 22, identified as wells 13, 17, 18 and 113 in the permit application, pursuant to section 14.2 of the Illinois Environmental Protection Act.

SPECIAL CONDITION 12. The use or operation of this facility shall be by or under the supervision of a Certified Class K operator.

SPECIAL CONDITION 13. The Outfall 002C Total Suspended Solids (TSS) loading rate shall be subtracted from the TSS loading rate determined for Outfall 002 when verifying compliance with the 18.6 lbs/day TSS load limit placed in the permit.

Additionally, when river water is used as a source of cooling water for Outfalls 002 and 004, a grab sample of the river water at the river water intake shall be taken and analyzed for TSS and the result shall be reported on the Discharge Monitoring Report (DMR) in mg/l. The background river TSS loading rate may be subtracted from the TSS loading rate for Outfalls 002 and 004 when verifying compliance with the TSS limits placed in the Effluent Limitations and Monitoring portion of the permit. If the TSS background river value is subtracted from the TSS results reported on the DMR for Outfalls 002 and/or 004, calculations shall be provided that demonstrate how the reported TSS values were determined.

#### Construction Authorization

Authorization is hereby granted to the above designee to construct the mine and mine refuse area described as follows:

The facility is an existing, approximately 802 acre inactive limestone quarry, aggregate processing area and wash plant, and cement processing plant, designated as the Dixon-Marquette Company, Dixon Plant, located in Sections 22, 27, 33 and 34, T22N, R9E of the 4th P.M. in Lee County, Illinois in Dixon. Facility operations include the crushing and stockpilling of limestone aggregate which is processed through a rotary kiln for the production of Portland Cement. Plant operation results in an average discharge of 0.06 MGD of noncontact cooling water from the powerhouse cooling towers from outfall 001A, 0.005 MGD of compressor cooling water from the Oid Clay Storage Building from outfall 001B, 0.009 MGD of raw mills and raw mill compressor cooling water from outfall 001C, 0.074 MGD of combined flows from outfalls 001A, 001B, 001C and storm water from outfall 001, 0.004 MGD of coal pile runoff from outfall 002A, 0.008 MGD of plant air compressors cooling water from cutfall 002B, 0.028 MGD of coal mill cooling water from outfall 002C, 0.132 MGD of precipitator compressors cooling water, preheater fans cooling water, analyzer water, Dracco baghouse fan cooling water, homogenizing silos compressors cooling water, #1,2,3 kiln piers discharge end cooling water from outfall 002D, 0.314 MGD of #1,2.3 finish mills cooling water. cement coolers, #1,2,3 finish mill compressor cooling water, #4 finish mill cooling water, #4 finish mill compressor cooling water. B pump compressors cooling water from outfall 002E, 0.48 MGD of combined flows from outfalls 002A-002E, and storm water from outfall 002. 0.008 MGD of packhouse compressors "A" pump cooling water from outfall 003A, 0.006 MGD of packhouse compressors instrument air cooling water from outfall 003B, 0.014 MGD of combined flows from outfalls 003A and 003B and storm water from outfall 003, 0.05 MGD of #1,2,3 kiln piers feed end cooling water, #4 kiln piers discharge end cooling water from outfall 004A, 0.10 MGD of #4 kiln compressors cooling water, wheelabrator baghouse fan motor bearings cooling water, #4 kiln piers feed end cooling water from outfall 004B, 0.15 MGD of combined flows from outfalls 004A and 004B and storm water runoff from an inactive limestone mining area and cement kiln dust landfill from outfall 004, storm water runoff on an intermittent basis from an inactive limestone mining area, aggregate wash plant sedimentation pond overflow and coal stock pile runoff from outfall 005, an intermittent discharge of storm water runoff from an inactive mining area and mine processing area from outfall 006, 0.229 MGD of storm water runoff from an inactive limestone mining area and cement kiln dust landfill from outfall 014 and 0.013 MGD of sewage treatment plant effluent from outfall 015. Sedimentation ponds will be constructed prior to outfalls 002A and 014. Outfalls 001, 002, 003, 005, 006, 014 and 015 and the flows tributary to these outfalls discharge to the Rock River. Outfall 004 and the flows tributary to outfall 004 discharge to an unnamed tributary of the Rock River.

The abandonment plan submitted with the application May 22, 1995 shall be executed and completed in accordance with Section 405.109 of Subtitle D: Mine Related Water Pollution.

This Authorization is issued subject to the following Special Condition(s). If such Special Conditions require additional or revised facilities, satisfactory engineering plan documents must be submitted to this Agency for review and approval.

If any statement or representation in the application is found to be incorrect, this permit may be revoked and the permittee thereupon waives all rights thereunder.

The issuance of this permit (a) shall not be considered as in any manner affecting the title of the premises upon which the mine or mine refuse area is to be located; (b) does not release the permittee from any liability for damage to person or property caused by or resulting from the installation, maintenance or operation of the proposed facilities; (c) does not take into consideration the structural stability of any units or parts of the project; and (d) does not release the permittee from compliance with other applicable statutes of the State of Illinois, or with applicable local laws, regulations or ordinances.

This permit may not be assigned or transferred. Any subsequent operator shall obtain a new permit from the Illinois Environmental Protection Agency.

There shall be no deviations from the approved plans and specifications unless revised plans, specifications and application shall first have been submitted to the Illinois Environmental Protection Agency and a supplemental permit issued.

The permit holder shall notify the Illinois Environmental Protection Agency (217/782-3637) immediately of an emergency at the mine or mine refuse area which causes or threatens to cause a sudden discharge of contaminants into the waters of Illinois and shall immediately undertake necessary corrective measures as required by Rule 405.111 under Chapter 1, Subtitle D: Mine Related Water Pollution of Illinois Pollution Control Board Rules and Regulations.

Final plans, specifications, application and supporting documents as submitted and approved shall constitute part of this permit and are identified in the records of the Illinois Environmental Protection Agency, by the permit number designated in the heading of this Section.

BY:DPG\98030407.DLK

#### Attachment H

#### Standard Conditions

#### Definitions

Act means the Illinois Environmental Protection Act, 415 ILCS 5 as Amended.

Agency means the Illinois Environmental Protection Agency.

Board means the Illinois Pollution Control Board.

Clean Water Act (formerly referred to as the Federal Water Pollution Control Act) means Pub. L 92-500, as amended. 33 U.S.C. 1251 et seq.

NPDES (National Pollutant Discharge Elimination System) means the national program for issuing, modfying, revoking and reissuing, terminating, monitoring and enforcing permits, and imposing and enforcing pretreatment requirements, under Sections 307, 402, 318 and 405 of the Clean Water Act.

USEPA means the United States Environmental Protection Agency.

Daily Discharge means the discharge of a pollutant measured during a calendar day or any 24-hour period that reasonably represents the calendar day for purposes of sampling. For pollutants with limitations expressed in units of mass, the "daily discharge" is calculated as the total mass of the pollutant discharged over the day. For pollutants with limitations expressed in other units of measurements, the "daily discharge" is calculated as the average measurement of the pollutant over the day.

Maximum Dally Discharge Limitation (daily maximum) means the highest allowable daily discharge.

Average Monthly Discharge Limitation (30 day average) means the highest allowable average of daily discharges over a celendar month, calculated as the sum of all delly discharges measured during a calendar month divided by the number of daily discharges measured during that month.

Average Weekly Discharge Limitation (7 day average) means the highest allowable average of daily discharges over a calendar week, calculated as the sum of all daily discharges measured during a calendar week divided by the number of daily discharges measured during that week.

Best Management Practices (BMPs) means schedules of activities, prohibitions of practices, maintenance procedures, and other management practices to prevent or reduce the pollution of waters of the State. BMPs also include treatment requirements, operating procedures, and practices to control plant site runoff, spillage or leaks, studge or waste disposal, or drainage from raw material storage.

Allquot means a sample of specified volume used to make up a total composite sample.

Grab Sample means an individual sample of at least 100 millilliers collected at a randomlyselected time over a period not exceeding 15 minutes.

24 Hour Composite Sample means a combination of at least 8 sample aliquots of at least 100 milliliters, collected at periodic intervals during the operating hours of a facility over a 24-hour period.

8 Hour Composite Sarriple means a combination of at least 3 sample aliquots of at least 100 milliliters, collected at periodic intervals during the operating hours of a facility over an 8-hour period.

Flow Proportional Composite Sample means a combination of sample aliquots of at least 100 milliliters collected at periodic intervals such that either the time interval between each aliquot or the volume of each aliquot is proportional to either the stream flow at the time of sampling or the total stream flow since the collection of the previous aliquot.

- (1) Duty to comply. The permittee must comply with all conditions of this permit. Any permit noncompliance constitutes a victation of the Act and is grounds for enforcement Ection, permit termination, revocation and reissuance, modification, or for deniel of a permit renewal application. The permittee shall comply with effluent standards or prohibitions established under Section 307(a) of the clean Water Act for toxic poliutants within the time provided in the regulations that establish these standards or prohibitions, even if the permit has not yet been modified to incorporate the requirement.
- (2) Duty to reapply. If the permittee wishes to continue an activity regulated by this permit after the expiration date of this permit, the permittee must apply for and obtain a new permit. If the permittee submits a proper application as required by the Agency no later than 180 days prior to the expiration date, this permit shall continue in full force and effect until the final Agency decision on the application has been made.
- (3) Need to halt or reduce activity not a defense, it shall not be a defense for a permittee in an enforcement action that it would have been necessary to half or reduce the permitted activity in order to maintain compliance with the conditions of this permit.
- (4) Duty to mitigate. The permittoe shall take all reasonable steps to minimize or prevent any discharge in violation of this permit which has a reasonable likelihood of adversely affecting human health or the environment.
- (5) Proper operation and maintenance. The permittee shall at all times properly operate and maintain all facilities and systems of treatment and control (and related appurtenances) which are installed or used by the permittee to achieve compliance with conditions of this permit. Proper operation and maintenance includes effective performance, adequate funding, adequate operator staffing and training, and adequate laboratory and process controls, including appropriate quality assurance procedures. This provision requires the operation of back-up, or auxiliary facilities, or similar systems only when necessary to schleve compliance with the conditions of the permit.

- 6) Permit actions. This permit may be modified, revoked and reissued, or termina for cause by the Agency pursuant to 40 CFR 122.62. The filing of a request by permittee for a permit modification, revocation and reissuance, or termination, or notification of planned changes or anticipated noncompliance, does not stay a permit condition.
- Property rights. This permit does not convey any property rights of any sort, or a exclusive privilege.
- (8) Duty to provide information. The permittee shall furnish to the Agency within reasonable time, any information which the Agency may request to determine whet cause exists for modifying, revoking and relissuing, or terminating this permit, of determine compliance with the permit. The permittee shall also furnish to the Agency upon request, copies of records required to be kept by this permit.
- (9) Inspection and ontry. The permittee shall allow an authorized representative of Agency, upon the presentation of credentials and other documents as may be requiby law, to:
  - Enter upon the permittee's premises where a regulated facility or activity located or conducted, or where records must be kept under the conditions of permit;
  - (b) Have access to and copy, at reasonable times, any records that must be k under the conditions of this permit;
  - (c) Inspect at reasonable times any facilities, equipment (including monitoring control equipment), practices, or operations regulated or required under permit; and
  - (d) Sample or monitor at reasonable times, for the purpose of assuring per compliance, or as otherwise authorized by the Act, any substances or parameter at any location.
- (10) Monitoring and records.
  - (a) Samples and measurements taken for the purpose of monitoring shall representative of the monitored activity.
  - (b) The permittee shall retain records of all monitoring information, including calibration and maintenance records, and all original strip chart recordings continuous monitoring instrumentation, copies of all reports required by permit, and records of all data used to complete the application for this permit, a period of all least 3 years from the date of this permit, measurement, repor application. This period may be extended by request of the Agency at any to
  - (c) Records of monitoring information shall include:
    - (1) The date, exact place, and time of sampling or measurements;
    - (2) The individual(s) who performed the sampling or measurements;
    - (3) The date(s) analyses were performed;
    - (4) The individual(s) who performed the analyses:
    - (5) The analytical techniques or methods used; and
    - (6) The results of such analyses.
  - (d) Monitoring must be conducted according to test procedures approved under CFR Part 136, unless other test procedures have been specified in this perr Where no test procedure under 40 CFR Part 136 has been approved, permittee must submit to the Agency a test method for approval. The permit shall calibrate and perform maintenance procedures on all monitoring a analytical instrumentation at intervals to ensure accuracy of measurements.
- (11) Signa bry requirement. All applications, reports or information submitted to: Agency shall be signed and certified.
  - (a) Application. All permit applications shall be signed as follows:
    - (1) For a corporation; by a principal executive officer of at least the level vice president or a person or position having ovarial responsibility environmental matters for the corporation;
    - (2) For a partnership or sole proprietorship; by a general partner or t proprietor, respectively; or
    - (3) For a municipality, State, Federal, or other public agency: by eithe principal executive officer or ranking elected official.
  - (b) Reports. All reports required by permits, or other information requested by the Agency shall be signed by a person described in paragraph (a) or by a disauthorized representative of that person. A person is a guly authorize representative only if;
    - (1) The authorization is made in writing by a person described in paragraph ( and)
    - (2) The authorization specifies either an individual or a position responsible the overall operation of the facility, from which the discharge originates, su as a plant manager, superintendent or person of equivalent responsibiliand
    - (3) The written authorization is submitted to the Agency.

- (c) Changes of Authorization. If an authorization under (b) is no longer accurate because a different individual or position has responsibility for the overall operation of the facility, a new authorization satisfying the requirements of (b) must be submitted to the Agency prior to or together with any reports, information, or applications to be signed by an authorized representative.
- (12) Reporting requirements.
  - (a) Planned changes. The permittee shall give notice to the Agency as soon as possible of any planned physical alterations or additions to the permitted facility.
  - (b) Anticipated noncompliance. The permittee shall give advance notice to the Agency of any planned changes in the permitted facility or activity which may result in noncompliance with permit requirements.
  - (c) Compliance schedules. Reports of compliance or noncompliance with, or any progress reports on, interim and final requirements contained in any compliance schedule of this permit shall be submitted no later than 14 days following each schedule date.
  - (d) Monitoring reports. Monitoring results shall be reported at the intervals specified elsewhere in this permit.
    - Monitoring results must be reported on a Discharge Monitoring Report (DMR).
    - (2) If the permittee monitors any pollutant more frequently than required by the permit, using test procedures approved under 40 CFR 136 or as specified in the permit, the results of this monitoring shall be included in the calculation and reporting of the data submitted in the DMR.
    - (3) Calculations for all limitations which require averaging of measurements shall utilize an arithmetic mean unless otherwise specified by the Agency in the permit.
  - (e) Twenty-four hour reporting. The permittee shall report any noncompliance which may endanger health or the environment. Any information shall be provided orally within 24 hours from the time the permittee becomes aware of the circumstances. A written submission shall also be provided within 5 days of the time the permittee becomes aware of the circumstances. The written submission shall contain a description of the noncompliance and its cause; the period of noncompliance, including exact dates and time; and if the noncompliance has not been corrected, the anticipated time it is expected to continue; and steps taken or planned to reduce, eliminate, and prevent reoccurrence of the noncompliance. The following shall be included as information which must be reported within 24 hours:
    - Any unanticipated bypass which exceeds any effluent limitation in the permit;
    - (2) Violation of a maximum daily discharge limitation for any of the pollutants listed by the Agency in the permit to be reported within 24 hours.

The Agency may waive the written report on a case-by-case basis if the oral report has been received within 24 hours.

- (f) Other noncompliance. The permittee shall report all instances of noncompliance not reported under paragraphs (12)(c), (d), or (e), at the time monitoring reports are submitted. The reports shall contain the information listed in paragraph (12)(e).
- (g) Other information. Where the permittee becomes aware that it failed to submit any relevant facts in a permit application, or submitted incorrect information in a permit application, or in any report to the Agency, it shall promptly submit such facts or information.
- (13) Transfer of permits. A permit may be automatically transferred to a new permittee if:
  - (a) The current parmities notifies the Agency at least 30 days in advance of the proposed transfer date;
  - (b) The notice includes a written agreement between the existing and new permittees containing a specific date for transfer of permit responsibility, coverage and liability between the current and new permittees; and
  - (c) The Agency does not notify the existing permittee and the proposed new permittee of its intent to modify or revoke and relissue the permit. If this notice is not received, the transfer is effective on the date specified in the agreement.
- (14) All manufacturing, commercial, mining, and silvicultural dischargers must notify the Agency as soon as they know or have reason to believe:
  - (a) That any activity has occurred or will occur which would result in the discharge of any toxic pollutent identified under Section 307 of the Clean Water Act which is not limited in the permit, if that discharge will exceed the highest of the following notification levels:
    - (1) One hundred micrograms per liter (100 ug/l);
    - (2) Two hundred inicrograms per liter (200 ug/l) for acroicin and acrylonkrite; five hundred micrograms per liter (500 ug/l) for 2,4-dinitrophenol and for 2methyl-4,6 dinitrophenol; and one milligram per liter (1 mg/l) for antimony.
    - (3) Five (5) times the maximum concentration value reported for that pollutant in the NPDES permit application; or

- (4) The level established by the Agency in this permit.
- (b) That they have begun or expect to begin to use or manufacture as an intermedia or final product or byproduct any toxic pollutant which was not imported a NPDES permit application.
- (15) All Publicly Owned Treatment Works (POTV/s) must provide adequate notice to I Agency of the following:
  - (a) Any new introduction of pollutants into that POTW from an indirect disci which would be subject to Sections 301 or 306 of the Clean Water Act if it directly discharging those pollutants; and
  - (b) Any substantial change in the volume or character of pollutants being introducing into that POTW by a source introducing pollutants into the POTW at the tilessuance of the permit.
  - (c) For purposes of this paragraph, adequate notice shall include information on the quality and quantity of efficient introduced into the POTW, and (ii) a anticipated impact of the change on the quantity or quality of efficient discharged from the POTW.
- (16) If the permit is issued to a publicly owned or publicly regulated treatment works, to permittee shall require any industrial user of such treatment works to comply will defer an equirements concerning;
  - (a) User charges pursuant to Section 204(b) of the Clean Water Act, and applied regulations appearing in 40 CFR 35;
  - (b) Toxic poliutant effluent standards and pretreatment standards pursuant to Sec 307 of the Clean Water Act; and
  - (c) Inspection, monitoring and entry pursuant to Section 308 of the Clean Water
- (17) If an applicable standard or limitation is promulgated under Section 301(b)(2)(C) at (D), 304(b)(2), or 307(a)(2) and that effluent standard or limitation is more strict than any effluent limitation in the permit, or controls a pollutant not limited appearsh, the permit shall be promptly modified or revoked, and reissued to contot that effluent standard or limitation.
- Any authorization to construct issued to the permittee pursuant to 35 il. Adm. 309.154 is hereby incorporated by reference as a condition of this permit.
- (19) The permittee shall not make any false statement, representation or certification in application, record, report, plan or other document submitted to the Agency or USEPA, or required to be maintained under this permit.
- (20) The Clean Water Act provides that any person who violates a permit con implementing Sections 301, 302, 308, 307, 308, 318, or 405 of the Clean Water is subject to a civil penalty not to exceed \$10,000 per day of such violation. A person who willfully or negligently violates permit conditions implementing Section 301, 302, 306, 307, or 308 of the Clean Water Act is subject to a fine of not lesses \$2,500 nor more than \$25,000 per day of violation, or by imprisonment for not than one year or both.
- (21) The Clean Water Act provides that any person who falstfies, tampers with knowingly renders inaccurate any monitoring device or method required to maintained under permit shall, upon conviction, be purished by a fine of not more \$10,000 per violation, or by imprisonment for not more than 6 months per violatio, by both.
- (22) The Clean Water Act provides that any person who knowingly makes any fall statement, representation, or certification in any record or other document submorrequired to be maintained under this permit shall, including monitoring reported to compliance or non-compliance shall, upon conviction, be punished by a of not more than \$10,000 per violation, or by imprisonment for not more than 6 monitoring per violation, or by both.
- (23) Collected screening, sturries, studget, and other solids shall be disposed of in a manner as to prevent entry of those wastes (or runoff from the wastes) into w of the State. The proper authorization for such disposal shall be obtained from Agency and is incorporated as part hereof by reference.
- (24) In case of conflict between these standard conditions and any other condition(s) shall govern.
- (25) The permittee shall comply with, in addition to the requirements of the permit, applicable provisions of 35 lii. Adm. Code, Subtitle C, Subtitle D, Subtitle E, and applicable orders of the Board.
- (26) The provisions of this permit are severable, and if any provision of this permit, a application of any provision of this permit is held invalid, the remaining provisions this permit shall continue in full force and effect.

(Rev. 3-13-98)

## VEGETATION ASSESSMENT REPORT SHORELINE AND BLUFF AREAS DIXON-MARQUETTE CEMENT KILN DUST LANDFILL

Prepared for:

Versar, Inc.

July 1999

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Mr. Gordon M. Stevens, P.G. Department Head, Geosciences Versar, Inc. Green Brook Executive Center 200 West 22nd Street, Suite 250 Lombard, IL 60148

22 July 1999

Re: Vegetation Assessment Report
Shoreline and Bluff Areas Adjoining

Shoreline and Bluff Areas Adjoining Dixon-Marquette Cement Kiln Dust Landfill Dixon, IL

Dear Mr. Stevens:

Buchanan Consulting Inc. (BCI) is pleased to submit this report as a deliverable from its services to Versar, Inc. to perform a vegetation assessment adjacent to the Dixon-Marquette Cement Kiln Dust (CKD) landfill located near Dixon, Illinois (Figure 1) as an indication if leachate from the CKD landfill operation is having an impact on the local terrestrial or shoreline ecosystem. The scope of work was completed in accordance with Schedule A of Versar's purchase order #07045 (corresponding to BCI's proposal dated 25 June 1999).

## **Understanding of the Project**

BCI's understanding of this project is that the CKD landfill at the Dixon-Marquette Cement Company facility is undergoing regulatory scrutiny. It has been deemed prudent and proactive to perform a vegetation assessment in the areas down-gradient from the CKD landfill to see if groundwater leachate and/or runoff from the CKD landfill is adversely impacting the bluff and shoreline environment adjacent to the Rock River. Assessing the vegetation adjacent to and down-gradient from the CKD landfill operation is being used as an environmental indicator because vegetation is a prominent component of the base of the ecological food chain, is dependent on and reflective of the immediate physical environment, and, therefore, will exhibit observable patterns in composition and form if influenced by contamination. The use of vegetation to indicate environmental contamination is a commonly used standard practice for property assessments (ASTM Standard E1527-94, "Standard Practice for Environmental Site Assessments: Phase 1 Environmental Site Assessment Process"). This assessment is voluntary and not regulatory driven.

BCI understands that its services and work product are confidential and are the property of Versar, Inc. for use by its client.

## Scope and Approach

BCI performed the following tasks to complete this assessment:

Collect Background Information - To ensure that the scope and approach was appropriate and accurate, BCI visited the site 23 June 1999 with Versar, Inc., met with the client, and performed a preliminary walkdown to identify appropriate locations for a representative vegetation assessment.

BCI reviewed literature on similar assessments of vegetation in river-side areas in Central Illinois and texts on regional flora to develop a context for what represents "normal" floristic composition. I/CI also collected information on the physical setting, soil types, disturbance history, etc. to ensure that comparisons between the observations made in the "potential impact zone" and the "control zone" are valid.

Perform Site Investigation - BCI visited the site 6 July 1999 and established four transects (i.e., linear observation routes through the representative vegetation communities), two transects in each of two types of vegetation communities (Figure 2). Within the area northwest of the CKD landfill, between the CKD landfill and the Rock River, are two primary vegetation types: the wooded bluff vegetation community and the river shoreline community. BCI established a transect in each of these two vegetation types within the potential impact zone adjacent to the CKD landfill and compared the observations to those taken from two similar transects in the control zone located north (up-gradient) outside the area of potential influence of the CKD landfil.

Per discussions with Versar (Mr. Gordon Stevens, personal communication, 23 June 1999), it is understood that groundwater travels toward the river through highly fractured dolomite/limestone, which forms the bluffs. The groundwater emerges at some points as active springs and also seeps into the river along the shoreline. A permitted surface water discharge also drains the landfill, traverses the bluff, and enters the river as a point discharge. If contamination is present in this groundwater and surface water at levels that could impact vegetation, these impacts would manifest themselves in the form of unusual vegetation species composition, or visible stress such as discoloration or odd growth forms.

BCI observed the vegetation along the designated transects and recorded the species composition of all existing strata (i.e., groundlayer, shrub, understory, and overstory).

These strata are defined as follows:

Groundlayer = all herbaceous (non-woody) species and seedlings of woody shrubs and trees growing on the ground;

Shrub stratum = woody shrubs, vines, and tree seedlings or young saplings growing above the groundlayer;

Understory = Woody species larger than 1" diameter but smaller than 4" diameter.

Overstory = Woody species larger than 4" diameter.

An indicator of abundance was recorded for each species in each stratum by ranking the species as: 1 - abundant and widespread; 2 - common; or 3 - infrequent or local. Qualitative observations of the condition of the vegetation were also recorded. Standard quantitative sampling methodologies were not employed for this assessment because the inherent variability of quantitative sampling methodologies renders them statistically invalid for comparing control and experimental vegetation transects in this manner. Qualitative observations by a trained plant ecologist have been adopted here as a more reliable means of detecting any abnormalities that might be due to operations at the subject facility.

The observations are summarized on data sheets for each stratum along each of the four transects. In addition, observations of wildlife or other evidence indicative of the health of the ecosystem were noted.

## Results

Setting - The assessment areas are located on the eastern shoreline and bluffs of the Rock River, just north of Dixon, Lee County, Illinois. This area forms the border of two principal geographical biotic divisions in Illinois: the Grand Prairie Division and Western Division (Jones 1963). The specific sampling transects are strongly influenced by very site-specific physical conditions that create a combination of soils, moisture regime and microclimate. In addition, the land traversed by the sampling transects has a history of human disturbance that also is manifested in the composition and distribution of the vegetation.

The two shoreline transects (Figures 3 and 4) were located between the Rock River waterway and the river frontage road (White Oak Road formerly Grand Detour Road) along a narrow strip of river bank. The soils are a combination of alluvial deposits and eroded silt-loams at the foot of the river bank grading into fractured limestone/dolomite and scree (loose rock pieces) overlain by a thin silt loam soil at the crest of the river bank. Slopes range from 15% in isolated areas where surface water outfalls from the bluffs have

deposited an alluvial delta to 80% in the most severely eroded river cuts. Generally, the banks slope 30% to 50%. Most of the length of the upper river bank in this area is occasionally mowed and maintained as roadside. This results in a component of the vegetation being reflective of early-successional disturbed sites close to the roadside. The lower river bank is occasionally subject to flooding and erosion/deposition. This influence is also reflected in certain components of the vegetation. A few scattered trees provide some shaded areas, especially along the up-gradient control transect, but most of the shoreline is in full sun and west-facing. At regular intervals are seeps and surface discharges from the fractured limestone/dolomite bluffs above and east of the river. The moisture regime, due to contributions from both the river and the seeps and springs, is wet to mesic.

The two wooded bluff transects (Figures 5 and 6) were located east of the river frontage road and beneath the crest of the bluff. The soils here are quite thin in most places, overlying fractured limestone/dolomite and scree near or at the surface in many places. Soil development is most advanced in level areas or depressed pockets within the bluffs, but generally the prevalent landform is steep scree slopes, 30% to 90%, and occasional bare limestone/dolomite escarpment. The predominance of limestone/dolomite contributes to a highly calcareous (calcium-rich) soil chemistry. The moisture regime varies from somewhat dry, as might be expected on a west-facing scree slope, to mesic in the vicinity of springs and seeps.

Vegetation Survey Results - The results of the qualitative surveys of each vegetation transect are presented in Tables 1 through 4 and can be summarized as follows:

## <u>Transect #1 Shoreline Transect Adjacent and Down-gradient from CKD Landfill Area</u>

Table 1 shows the results for this transect. The most prominent species in the ground layer are a mix of typical central Illinois floodplain species (orange jewelweed and Virginia creeper) and weedy invasive and non-native species reflective of physical disturbances (reed canarygrass and black mustard). A total of 30 groundlayer species were observed, indicating a reasonably diverse vegetation matrix. In the shrub stratum, the cominant species (indigo bush, Virginia creeper, and riverbank grape) are all typical native Illinois floodplain and riparian residents. A total of 16 species of shrubs and tree seedlings were observed, almost all of which were typical native species for this situation. Only a few large scattered trees occur along the shoreline, consisting of white ash and cottonwood. Again, these are typical of this type of vegetation community. Per Swink and Wilhelm (1979) Plants of the Chicago Region, the association of species found here are generally reflective of other similar plant associations found elsewhere in north-central and northeast Illinois.

All of the vegetation strata were vigorous, healthy, and showed no signs of stress, such as discoloration, malformation, or necrosis.

## Transect #2 Shoreline Transect Up-gradient from CKD Landfill Area

This transect was selected as a background or control area outside of the influence of the CKD landfill area. Table 2 presents the observations from this transect. To make the comparison of the vegetation of this transect with that of the down-gradient transect, it is important to note that there were a few significant physical differences between transects. First, the up-gradient transect is shorter in length, due to the fact that its northern end abuts cropland and a camping/recreational area so that sampling beyond that point was untenable. Second, portions of the up-gradient transect are shaded by trees more than the down-gradient transect, and the ground layer composition is influenced somewhat by this shade.

Even so, the composition of the up-gradient transect is quite similar to the down-gradient transect. The two dominant species in the groundlayer are orange jewelweed and reed canarygrass, similar to the down-gradient composition. Species diversity was almost identical, with 29 groundlayer species recorded, of which 24 were also recorded in the down-gradient transect. Of the five species that were found in the up-gradient transect not found in the down-gradient, four are shade-loving. The shrub stratum also matches quite closely between the up-gradient transect and down-gradient transect. Both transects include Virginia creeper and riverbank grape as dominants. The notable difference between the two transects in shrub composition is that indigo bush is prominent in the down-gradient transect and lacking in the up-gradient. Indigo bush is not particularly shade tolerant, which might explain its absence in the up-gradient transect. This does not indicate that the down-gradient transect is relatively impacted or degraded, in fact, to the contrary, indigo bush is a fairly conservative species (Quality Rating of 6 by Swink and Wilhelm 1979) which means that it is fairly intolerant of disturbance. In a similar vein, the existence in both transects of glade mallow, a very high quality species intolerant of disturbance, underscores this point.

Again, the vegetation in the up-gradient transect was healthy, vigorous, and showed no signs of adverse environmental impacts.

# Transect #3 Bluff Woods Transect Adjacent and Down-gradient from CKD Landfill Area

The observations of species in the various strata in this survey transect are presented in Table 3. The groundlayer was quite diverse, with white snakeroot, Virginia creeper, and poison ivy the most prevalent species. A total of 32 species were observed in the groundlayer, indicating a fairly high diversity. Almost all of these species are native and typical of a calcareous river bluff woodland. The shrub stratum was also fairly diverse,

with honeysuckle and choke cherry dominant, and a total of 12 species of tree reproduction and shrubs observed. The understory was composed of 10 species of tree saplings and understory trees, of which hackberry was most widespread. The overstory was dominated by boxelder and contained a total of 14 tree species; however, both the understory and overstory had a fairly even mix of tree species, indicating good balance and healthy reproduction. Nearly all of the species, with the exception of white mulberry which is a non-native adventive species, are typical native residents of river-side slopes in this part of Illinois. All strata appeared healthy and vigorous.

### Transect #4 Bluff Woods Transect Up-gradient from CKD Landfill Area

Table 4 presents the observations of species in this transect. Again, for comparative purposes, this transect was not quite as long as Transect #3 adjoining the CKD landfill area due to changes in topography that limited the survey length. However, the results indicate the composition and structure of the two transects are similar and, therefore, underscore the validity of the comparisons. In fact, all of the 25 species observed in the groundlayer of the up-gradient transect, also appear in the adjoining transect. The shrub stratum was dominated by honeysuckle, as was the adjoining transect, and in general, all strata were similarly composed. The slightly lower diversity in the up-gradient transect probably simply reflects the smaller sampling area due to the aforementioned topographic limitations on the transect placement. All strata appeared healthy and vigorous.

Discussion and Conclusions - In general, the vegetation both within and outside of the potential zone of influence of the CKD landfill operation appear quite similar and in excellent condition. No evidence of impacts from the CKD landfill operation was observed, and no discernible differences in composition, form, or health were observed between the up-gradient transects and those adjoining or down-gradient from the CKD landfill.

The shoreline transects both contain a mix of plant species reflective of some amount of physical disturbance, which is to be expected in this setting with the periodic flooding and road construction/maintenance. The wooded bluff transects reflect a fairly high quality river-side woods in all strata. All transects support a species composition and structure that are quite typical of this location and physical setting. When comparing the results of the survey to other similar reports from the Illinois River valley (Buchanan 1975), the Sangamon River basin (Jones and Bell 1974), and the typical associations reported by Swink and Wilhelm (1979) for the north-central and northeast Illinois region, this conclusion is strongly supported.

The differences that exist between the up-gradient sampling transects and the adjoining or down-gradient transects are easily attributable to microclimate, topography and natural diversity. Many of the species found in both transects, but most notably in the transects

adjoining or down-gradient from the landfill operation, are conservative species intolerant to disturbance. The bluff woods in particular are fairly high quality in terms of vegetation. While a few invasive species such as common buckthorn and white mulberry occur, they are not dominant as in other degraded woodlands throughout the region. Relatively few exotic non-native species were observed in any transect. Numerous species observed in both the shoreline and bluff woods survey transects are assigned quality ratings by Swink and Wilhelm of 5 or higher (on a scale of 0 - 10). If these shoreline and woodland communities were being significantly and adversely impacted by the landfill operations, such conservative species would probably not occur.

Corollary observations were made of wildlife usage as an indication of ecosystem health. While certainly not exhaustive or complete, the presence of such higher level consumers as hawks and turtles, the abundance of frogs, fish hatchlings in the outfall points of the springs from the bluffs, and nesting birds are suggestive of a healthy functioning ecosystem.

Thank you very much for the opportunity to serve you. Please contact BCI with any questions or comments.

Respectfully submitted,
Warren J. Buchanan

Warren J. Buchanan

President

### **References Cited**

- 1. Buchanan, W. J. 1975. Vegetation monitoring. <u>IN Terrestrial monitoring program for</u> the construction phase of the LaSalle County Station 1975. Second Annual Report to Commonwealth Edison Company, Chicago. Chapter 1. 66 pages.
- Jones, A. G. and D. T. Bell. 1974. Vascular plants of the Sangamon River Basin: annotated checklist and ecological summary. Univ. Il. College of Ag. Ag. Expt. Sta. Bull. 746. 77 pages.
- 3. Jones, G. N. 1963. Flora of Illinois. Third ed. American Midland Naturalist: Monograph No. 7. 401 pages.
- 4. Swink, F. and G. Wilhelm. 1979. Plants of the Chicago Region. Third ed. The Morton Arboretum, Lisle, IL. 922 pages.

**FIGURES** 

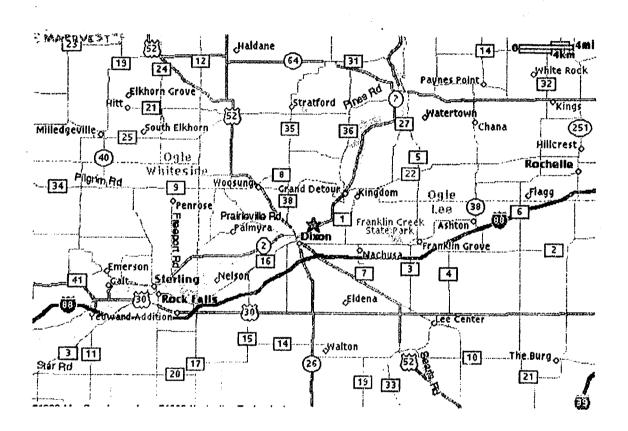


Figure 1. Site location

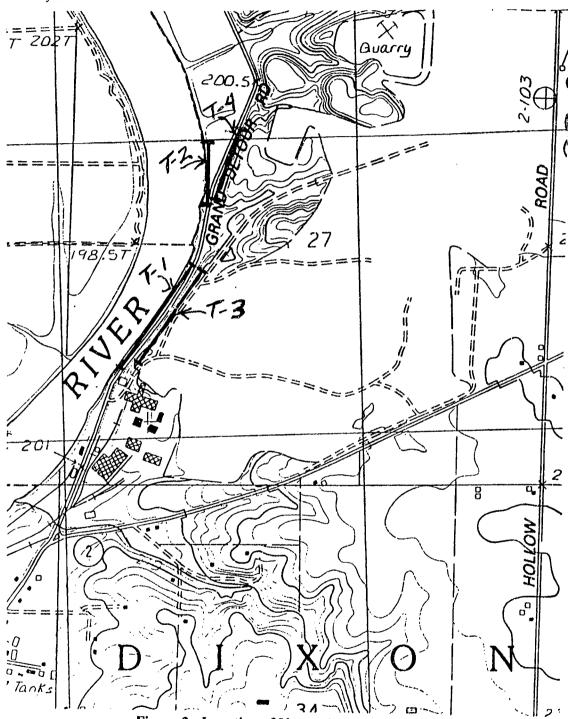


Figure 2. Location of Vegetation Transects

Figure 3. Shoreline Vegetation Transect #1
Adjacent to and Down-Gradient from CKD Landfill

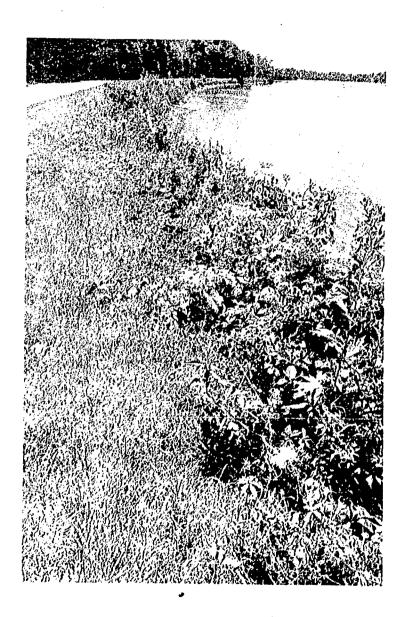
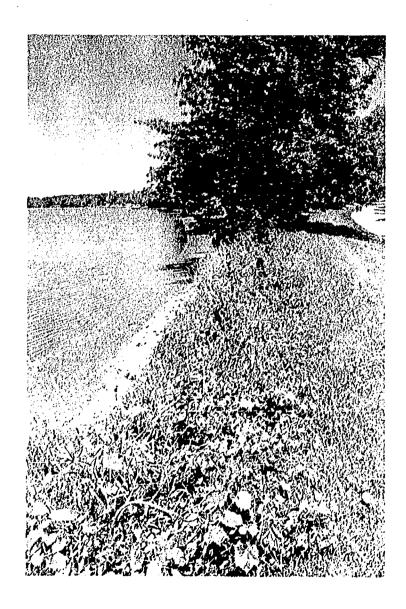


Figure 4. Shoreline Vegetation Transect #2
Up-Gradient from CKD Landfill



## Buchanan Consulting, Inc.

Figure 5. Bluff Woods Vegetation Transect #3
Adjacent to and Down-Gradient from CKD Landfill



## Buchanan Consulting, Inc.

Figure 6. Bluff Woods Vegetation Transect #4
Up-Gradient from CKD Landfill



## Buchanan Consulting, Inc.

Mr. Gordon M. Stevens, P.G.: Versar, Inc. Vegetation Assessment, Dixon-Marquette Cement, Dixon, IL 22 July 1999

**TABLES** 

## Table 1. List of Species and Abundance Index for Transect #1 Shoreline Transect Adjacent to Landfill Area

('Note: 1=abundant, widespread; 2=common; 3=infrequent, local)

## Groundlayer

Scientific Name	Common Name	Abundance*
Brassica nigra	Black mustard	l
Impatiens capensis	Orange jewelweed	1
Parthenocissus quinquefolia	Virginia creeper	1
Phalaris arundinacea	Reed canarygrass	1
Bromus inermis	Smooth brome	2
Convolvulus sepium	Hedge bindweed	2
Erigeron philadelphicus	Marsh fleabane	2
Heliopsis helianthoides	False sunflower	2
Napaea dioica	Glade mallow	2
Nepeta cataria	Catnip	2
Oenothera biennis	Common evening primrose	2
Rhus radicans	Poison ivy	2
Rudbeckia laciniata	Wild golden glow	2
Silphium perfoliatum	Cup plant	2
Urtica gracilis	Tall nettle	2
V <sup>rei</sup> s riparia	Riverbank grape	2 2 2 3 3
A, Jeynum cannabinum	Dogbane	3
Asclepias incarnata	Marsh milkweed	3
Asclepias syriaca	Common milkweed	3
Aster sp.	Aster species	3
Cicuta maculata	Water hemlock	3
Eupatorium rugosum	White snakeroot	3
Heracleum maximum	Cow parsnip	3
Ipomoea purpurea	Common morning glory	3
Menispermum canadense	Moonseed	3
Polygonatum biflorum	Solomon's scal	3
Smilax tamnoides hispida	Bristly catbrier	3
Solidago altissima	Tall goldenrod	3
Teucrium canadense	Germander or wood sage	3
Thalictrum polygamum	Meadow rue	3

## Shrub Stratum

Scientific Name	Common Name	Abundance
Amorpha fruticosa	Indigo bush	1
Parthenocissus quinquefolia	Virginia creepor	1

# Table 1. (continued) List of Species and Abundance Index for Transect #1 Shoreline Transect Adjacent to Landfill Area

('Note: 1=abundant, widespread; 2=common; 3=infrequent, local)

## Shrub Stratum (continued)

Scientific Name	Common Name	Abun lance
Vitis riparia	Riverbank grape	1
Acer negundo	Boxelder	2
Acer saccharinum	Silver maple	2
Fraxinus americana	White ash	2
Rhus radicans	Poison ivy	2
Celtis occidentalis	Hackberry	3
Cornus racemosa	Gray dogwood	3
Lonicera prolifera	Honeysuckle	3
Morus alba	White mulberry	3
Rosa multiflora	Multiflora rose	3
Prunus virginiana	Choke cherry	3
Sambucus canadensis	Elderberry	3
Tilia americana	Basswood	3
Ulmus rubra	Slippery elm	3

## Tree Stratum (few scattered trees)

Scientific Name	Common Name	Abundance
Fraxinus americana	White ash	3
Populus deltoides	Cottonwood	3

## Table 2. List of Species and Abundance Index for Transect #2 Shoreline Transect Upgradient from Landfill Area

('Note: 1=abundant, widespread; 2=common; 3=infrequent, local)

## Groundlayer

Scientific Name	Common Name	Abundance '
Impatiens capensis	Orange jewelweed	1
Phalaris arundinacea	Reed canarygrass	1
Brassica nigra	Black mustard	2
Bromus inermis	Smooth brome	2
Convolvulus sepium	Hedge bindweed	2
Heliopsis helianthoides	False sunflower	2
Napaea dioica	Glade mallow	2
Parthenocissus quinquefolia	Virginia creeper	2
Rhus radicans	Poison ivy	2
Rudbeckia laciniata	Wild golden glow	2
Silphium perfoliatum	Cup plant	2
Teucrium canadense	Germander or wood sage	2
Vitis riparia	Riverbank grape	2
Actinomeris alternifolia	Wingstem	3
Asclepias incarnata	Marsh milkweed	3
Asclepias syriaca	Common milkweed	3
Aster sp.	Aster species	3
Elymus virginicus	Virginia wildrye	3
Heracleum maximum	Cow parsnip	3
Laportea canadensis	Wood nettle	3
Menispermum canadense	Moonseed	3
Oenothera biennis	Common evening primrose	3
Polygonatum biflorum	Solomon's seal	3
Sanguinaria canadensis	Bloodroot	3
Silene cucubalus	Bladder campion	3
Smilax tamnoides hispida	Bristly catbrier	3
Solidago altissima	Tall goldenrod	3
Thalictrum polygamum	Meadow rue	3
Urtica gracilis	Tall nettle	3

## **Shrub Stratum**

Scientific Name	Common Name	<u>Abundance</u>
Parthenocissus quinquefolia	Virginia creeper	1
Vitis riparia	Riverbank grape	1
Rhus radicans	Poison ivy	2

## Table 2. (continued) List of Species and Abundance Index for Transect #2 Shoreline Transect Upgradient from Landfill Area

('Note: 1=abundant, widespread; 2=common; 3=infrequent, local)

## **Shrub Stratum (continued)**

Scientific Name	Common Name	Abundance'
Celtis occidentalis	Hackberry	3
Corris racemosa	Gray dogwood	3
Lonicera prolifera	Honeysuckle	3
Morus alba	White mulberry	3
Rosa multiflora	Multiflora rose	3
Sambucus canadensis	Elderberry	3
Staphylea trifolia	Bladdernut	3
Tilia americana	Basswood	3
Ulmus rubra	Slippery elm	3
Viburnum opulus	Flighbush cranberry	. 3

## Tree Stratum (few scattered trees)

Scientific Name	Common Name	<b>Abundance</b>
Acer negundo	Boxelder	3
Acer saccharinum	Silver maple	3
Fraxinus americana	White ash	3
Ulmus americana	American elm	3

# Table 3. List of Species and Abundance Index for Transect #3 Bluff Woods Transect Adjacent to Landfill Area

('Note: 1=abundant, widespread; 2=common; 3=infrequent, local)

## Groundlayer

Scientific Name	Common Name	Abundance*
Eupatorium rugosum	White snakeroot	1
Parthenocissus quinquefolia	Virginia creeper	1
Rhus radicans	Poison ivy	1
Alliaria officinalis	Garlic mustard	2
Aster shortii	Short's aster	2
Campanula americana	Tall beliflower	2
Carex laxiflora	Wood sedge.	2
Hepatica acutiloba	Sharp-lobed hepatica	2
Sanguinaria canadensis	Bloodroot	2
Sanicula gregaria	Black snakeroot	2
Smilacina racemosa	False Solomon's seal	2
Smilax tamnoides hispida	Bristly catbrier	2
Anemone virginiana	Thimble flower	3
Aquilegia canadensis	Columbine	3
Arabis laevigata	Smooth bank cress	3
Celtis occidentalis	Hackberry	3
Circaea quadrisulcata	Enchanter's nightshade	3
Desmodium glutinosum	Pointed tick-trefoil	3
Elymus villosus	Silky wildrye	3
Galium concinnum	Shining bedstraw	3
Geum canadense	White avens	3
Helianthus strumosus	Woodland sunflower	3
Hydrophyllum virginianum	Virginia waterleaf	3
Impatiens capensis	Orange jewelweed	3
Menispermum canadense	Moonseed	3
Osmorhiza claytoni	Sweet cicely	3
Physalis heterophylla	Clammy ground cherry	3
Polygonatum biflora	Solomon's seal	3
Prenanthes alba	Lion's foot	3
Sanguinaria canadensis	Bloodroot	3
Solidago flexicaulis	Zigzag goldenrod	3
Trillium recurvatum	Prairie wake robin	3

# Table 3. (continued) List of Species and Abundance Index for Transect #3 Bluff Woods Transect Upgradient from Landfill Area

('Note: 1=abundant, widespread; 2=common; 3=infrequent, local)

## **Shrub Stratum (continued)**

Scientific Name	Common Name	Abundance*
Vitis riparia	Riverbank grape	2
Celtis occidentalis	Hackberry	3
Fraxinus quadrangulata	Blue ash	3
Physocarpus opulifolius	Ninebark	3

## **Understory Stratum**

Scientific Name	Common Name	Abundance
Celtis occidentalis	Hackberry	1
Fraxinus americana	White ash	2
Fraxinus quadrangulata	Blue ash	2
Acer negundo	Boxelder	3
Acer saccharum	. Sugar maple	3
Juglans nigra	Black wainut	3
Ostrya virginiana	Hop hornbeam	3
Ulmus rubra	Slippery elm	3

## **Overstory Stratum**

Scientific Name	Common Name	<u>Abundance</u>
Celtis occidentalis	Hackberry	2
Fraxinus quadrangulata	Blue ash	2
Quercus macrocarpa	Burr oak	2
Tilia americana	Basswood	2
Fraxinus americana	White ash	3
Juglans nigra	Black walnut	3
Populus deltoides	Cottonwood	3
Ulmus rubra	Slippery elm	3

## Table 4. List of Species and Abundance Index for Transect #4 Bluff Woods Transect Upgradient from Landfill Area

('Note: 1=abundant, widespread; 2=common; 3=infrequent, local)

#### Groundlayer

	Scientific Name	Common Name	Abundance'
	Parthenocissus quinquefolia	Virginia creeper	1
	Aster shortii	Short's aster	2
	Eupatorium rugosum	White snakeroot	2
	Osmorhiza claytoni	Sweet cicely	2
	Prenanthes alba	Lion's foot	2
	Rhus radicans	Poison ivy	$\overline{2}$
	Sanguinaria canadensis	Bloodroot	2
	Sanicula gregaria	Black snakeroot	2
	Alliaria officinalis	Garlic mustard	3
	Anemone virginiana	Thimble flower	3
P	Aquilegia canadensis	Columbine •	3
	Arabis laevigata	Smooth bank cress	3
	Campanula americana	Tall bellflower	3
	Carex laxiflora	Wood sedge.	3
	Celtis occidentalis	Hackberry	3
	Elymus villosus	Silky wildrye	3
	Galium concinnum	Shining bedstraw	3
	Geum canadense	White avens	3
	Hydrophyllum virginianum	Virginia waterleaf	3
	Impatiens capensis	Orange jewelweed	3
	Menispermum canadense	Moonseed	3
	Smilacina racemosa	False Solomon's seal	3
	Smilax tamnoides hispida	Bristly catbrier	3
	Solidago flexicaulis	Zigzag goldenrod	3
	Trillium recurvatum	Prairie wake robin	3

## Shrub Stratum

Scientific Name	Common Name	Abundance
Lonicera prolifera	Honeysuckle	1
Fraxinus americana	White ash	2
Prunus virginiana	Choke cherry	2
Rhamnus cathartica	Common buckthorn	2
Ribes missouriense	Missouri gooseberry	$\overline{2}$
Smilax tamnoides hispida	Bristly catbrier	2
Staphylea trifolia	Bladdernut	2

# Table 4. (continued) List of Species and Abundance Index for Transect #4 Bluff Woods Transect Adjacent to Landfill Area

(Note: 1=abundant, widespread; 2=common; 3=infrequent, local)

#### **Shrub Stratum**

<u>Scientific Name</u>	Common Name	<u>Abundance</u>
Lonicera prolifera	Honeysuckle	1
Prunus virginiana	Choke cherry	i
Rhamnus cathartica	Common buckthorn	2
Rhus radicans	Poison ivy	2
Smilax tamnoides hispida	Bristly catbrier	2
Staphylea trifolia	Bladdernut	2
Vitis riparia	Riverbank grape	2
Acer negundo	Boxelder	3
Juglans nigra	Black walnut	3
Physocarpus opulifolius	Ninebark	3
Rhus glabra	Smooth sumac	3
Ribes missouriense	Missouri gooseberry	3

## **Understory Stratum**

Scientific Name	Common Name	Abundance
Celtis occidentalis	Hackberry	1
Acer negundo	Boxelder	2
Fraxinus americana	White ash	2
Ostrya virginiana	Hop hornbeam	2
Staphylea trifoliata	Bladdernut	2
Acer saccharum	Sugar maple	3
Fraxinus quadrangulata	Blue ash	3
Juglans nigra	Black walnut	3
Morus alba	White mulberry	3
Quercus macrocarpa	Burr oak	3

## **Overstory Stratum**

Scientific Name	Common Name	Abundance
Acer negundo	Boxelder	
Celtis occidentalis	Hackberry	2
Fraxinus americana	White ash	2
Fraxinus quadrangulata	Blue ash	2

## Table 4. (continued) List of Species and Abundance Index for Transect #4 Bluff Woods Transect Adjacent to Landfill Area

('Note: 1=abundant, widespread; 2=common; 3=infrequent, local)

## **Overstory Stratum (continued)**

Scientific Name	Common Name	<b>Abundance</b>
Quercus borealis	Red oak	<b>2</b>
Quercus macrocarpa	Burr oak	2
Acer saccharum	Sugar maple	3
Juglans nigra	Black walnut	3
Morus alba	White mulberry	3
Morus rubra	Red mulberry	3
Ostrya virginiana	Hop hornbeam	<b>3</b>
Populus deltoides	Cottonwood	3
Tilia americana	Basswood	3
Ulmus rubra	Slippery elm	. 3

## **DETECTION MONITORING PROGRAM**

for Dixon Marquette Cement Company Dixon, Illinois

January 2001

Prepared by

Preston Engineering, Inc. 4436 North Brady Street Davenport, IA 52806 319/388-8288

#### 1.0 INTRODUCTION

During March and April 1998, nine monitoring wells (MW1-S, MW1-D, MW2-S, MW2-D, MW2-M, MW3-S, MW3-D, MW4-S, and MW4-D) were installed around the CKD disposal site at the Dixon Marquette Cement plant in Dixon, Illinois. The wells were installed in pairs at three sites, and three wells were installed at one site. Each well was assigned a number designating which well cluster it was in, and a letter designating the location of the screened interval. Letter designations were S and D, which designated screened intervals at the base of the Pecatonica Formation and the top of the St. Peter Formation, respectively. One well, completed in the Glenwood Formation, was designated MW2-M, and was used solely for conducting hydraulic conductivity testing. Each well, except for MW2-M, has a 10 foot long, 2 inch diameter screen at its base.

In May 1998 two leachate wells were installed in the CKD disposal area. The wells were designated LW-1 and LW-2. The base of the screen in both wells was positioned at the contact between CKD and the underlying Pecatonica Formation. One of the wells, LW-2, was dry and was later abandoned. Well LW-1 has consistently had several feet of water in it. The water in the well is assumed to have been in contact with the surrounding CKD and is considered to be leachate. An additional leachate well, LW-3, was installed in the CKD disposal area in July 1999. No ground water was noted in the CKD during drilling. Groundwater was found in the uppermost part of the Pecatonica Formation, which was directly under the CKD.

In June 1999 four pairs of wells were installed along White Oak Lane on the west side of Dixon Marquette property. In each pair, one well is screened across the water table, and one is screened at the top of the St. Peter Formation. The wells were designated MW5-S, MW5-D, MW6-S, MW6-D, MW7-S, MW7-D, MW8-S, and MW8-D.

In June 1999 a pair of wells was installed south of the CKD disposal area between wells MW2 and MW3. They were designated MW9-S and MW9-D. The screen in well MW9-S was placed at the base of the Pecatonica Formation, and the screen in MW9-D was placed at the top of the St. Peter Formation.

In July 1999 wells were added that were screened across the potentiometric surface, or as close to it as practical, at five locations. These wells were designated MW1-WT, MW2-WT, MW3-WT, MW4-WT, and MW9-WT.

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Seven piezometers (P1 through P7) were added southeast, east, and north of the landfill. They are narrow diameter PVC pipes that run from just above the surface to about 15 feet deep. They are used to measure the potentiometric surface near the top of the uppermost aquifer, and they are not included in the sampling program. Figure 1 is a map of the site with the well and piezometer locations shown, and Table 1 lists well data.

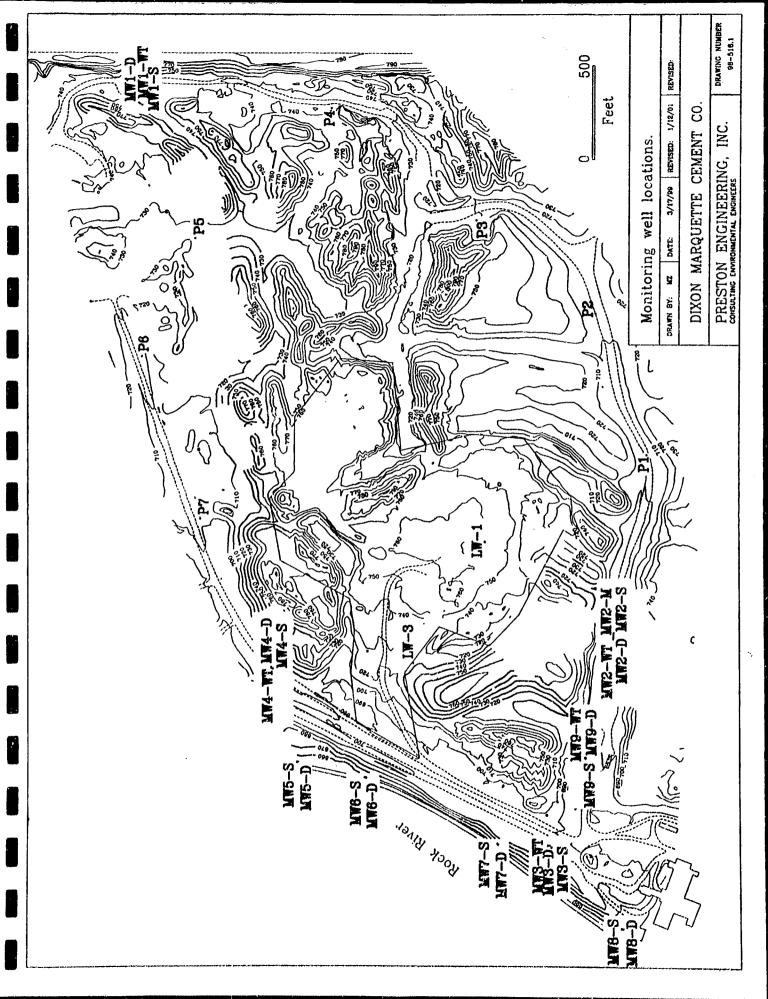


Table 1. Monitoring well data.

Well	Total Depth (ft from top of casing)	Formation screened	
MW1-S	42.05	Pecatonica	
MW1-D	80.5	St. Peter	
MWI-WT	15.5	Pecatonica	
MW2-S	41.85	Pecatonica	
MW2-D	62.25	St. Peter	
MW2-M	46.20	Glenwood	
MW2-WT	15.5	Pecatonica	
MW3·S	61.5	Pecatonica	
MW3-U	80.6	St. Peter	
MW3-WT	37.5	Pecatonica	
MW4-S	44.1	Pecatonica .	
MW4-D	63.2	St. Peter	
MW4-WT	25	Pecatonica	
MW5-S	15	Alluvium*	
MW5-D	47.5	St. Peter	
MW6-S	15	Alluvicm*	
MW6-D	40	St. Peter	
MW7-S	15	Pecatonica	
MW7-D	34.5	Si. Peter	
MW8-S	18	Pecatonica	
MW8-D	40	St. Peter	
MW9-S	37.5	Pecatonica	
MW9-D	59.5	St. Peter	
MW9-WT	15.5	Pecatonica	
LW-1	65	CKD fill	
LW-3	44.5	Pecatonica	
Spring	not applicable	emanates from Pecatonica	

<sup>\*</sup>The Pecatonica Formation has been croded at this location and replaced with alluvium of the Rock River.

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The groundwater in and around the CKD disposal area will be monitored quarterly. The sampling plan is designed to detect changes in groundwater quality that result from contamination by leachate derived from the landfill.

Groundwater flow beneath the landfill is roughly east to west, or toward the Rock River. One well triplet (MW1-S, MW1-D, and MW1-WT) was placed on the east, or upgradient side of the landfill. These wells monitor the groundwater quality in an area that has not been affected by CKD disposal. Data from the upgradient wells will be used in comparisons with data from other wells to determine if leachate has affected the groundwater quality.

Wells at MW2, MW3, MW4, and MW9 are used to detect leachate in the groundwater in areas close to the CKD and downgradient from it. Well pairs at MW5, MW6, MW7, and MW8 are used to determine if leachate has affected the groundwater near the Rock River, and to determine if the groundwater quality may affect the surface water quality in the river.

Two wells, LW-1 and LW-3, are used to characterize groundwater beneath the landfill.

Groundwater in these wells may have been in contact with the CKD. Data from these wells will be used to determine which parameters should be monitored in other wells.

Water elevations in all of the wells and piezometers will be measured before each sampling event. The data can be used to create potentiometric maps of the groundwater. Maps will be made from data from groups of wells that are screened at the same depth in the same formation.

#### 2.0 SAMPLING METHODOLOGY

Prior to field activities, arrangements will be made with a commercial laboratory that is certified in the State of Illinois. The lab will provide sample containers and will place the proper preservative in each container prior to shipping. A list of the container types and preservatives is presented in Table 2.

The following information from each well will be recorded at the time of sampling:

Date and time of sampling

Depth to water from the top of the casing prior to purging

The volume of water that was purged prior to sampling

The appearance of the water

The temperature, pH, conductivity, and turbidity of the groundwater at the time of sampling

Wells MW1-S, MW1-D, MW1-WT, MW2-S, MW2-D, MW2-WT, MW3-S, MW3-D, MW3-WT, MW4-S, MW4-D, MW4-WT, MW5-S, MW5-D, MW6-S, MW6-D, MW7-S, MW7-D, MW8-S, MW8-D, MW9-S, MW9-D, MW9-WT, LW-1, LW-3, the Rock River (two locations), and the spring that is west of MW-3 will be sampled. We'ls will be purged prior to sampling. Sampling will take place within 24 hours of purging. The water will be field tested for turbidity. If the turbidity is 5 NTU or less, samples will not be field filtered for metals. If the turbidity exceeds 5 NTU, two samples for metals will be collected: one that is unfiltered, and one that is field filtered through a filter with a pore size of 0.45 µm or less prior to preservation. All of the samples will be kept on ice in the field and during delivery to the lab. Samples will be packed in coolers with ice, sealed, and delivered to the laboratory by overnight courier. A chain of custody form will a company the shipment. An example chain of custody form is attached. This form, or an equivalent form supplied by the laboratory, will be used.

Sampling of the Rock River will take place quarterly. Sampling will consist of collection of river water from a location upgradient from the landfill (about 1000 feet upriver from MW5) and from a location downgradient from the landfill (near MW8). Sampling will take place from the shore on the left bank. The river water will be analyzed for the same parameters as the groundwater. Samples will first be collected from the upstream location, then the samples from the downstream location will be collected. Field measurements of temperature, pH, and specific conductivity will be made. The river samples for dissolved iron will be passed through a filter with a pore size of 0.45  $\mu$ m or less prior to preservation.

In addition to monitoring for compounds that exceed the background concentrations, the list of organic chemicals in 40 CFR 141.40, and any other organic chemical for which a groundwater quality standard or criterion has been adopted pursuant to Section 14.4 of the Act or Section 8 of the Illinois Groundwater Protection Act, will be analyzed biennially.

Table 2. Chemical Parameters to be Analyzed.

Parameter	Method	Container	Preservative	Holding Time
Volatile organics	8260	2 40 ml glass vials	HCl, 4°C	14 days
Metals	200.8	125 ml HDPE	HNO3, 4°C	6 months
pН	4500	250 ml HDPE	4°C	immediate
Calcium, Potassium, Sodium, Chloride, Fluoride, Nitrate, Sulfate, Alkalinity, Ammonia, Dissolved solids, TOC	200.8, 4110, 2320, 2540C, 4500, 5310	250 ml HDPE	4°C	14 days

#### 3.0 ANALYTES

Quarterly samples from all 28 locations will be analyzed for the following. Practical quantitation limits will be those specified by USEPA SW846 laboratory methods.

#### Benzene

- 1.1-dichloroethane
- 1,1-dichloroethene
- 1,1,1-trichloroethane

Toluene

Aluminum

Antimony

Arsenic

**Barium** 

Beryllium

Boron

Cadmium

Calcium

Chromium

Cobalt

Copper

Iron

Lead

Magnesium

Manganese

Mercury

Nickel

Potassium

Selenium

Sodium

Thallium

Vanadium

Zinc

Ammonia nitrogen

Chloride

COD

Fluoride

Nitrate as N

Sulfate

Alkalinity, phenol.

Alkalinity, bicarbonate

Alkalinity, total

Total organic carbon

Total dissolved solids

pН

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These parameters have been previously detected in downgradient wells at concentrations exceeding the background concentrations.

#### 4.0 QUALITY CONTROL

Additional samples for quality control will be submitted to the laboratory. Split samples will be collected from wells MW1-S and MW3-S (both screened in the Pecatonica Formation), and MW7-S. The samples will be labeled as split samples, but no well numbers will identify them. The split samples will be analyzed for the same parameters as the other samples from the same wells. These samples will be used as a check on sampling procedures and reproducibility at the laboratory.

A trip blank consisting of pure water will be prepared by the laboratory and analyzed for organic compounds by Method 8260 and metals by Method 200.8. The trip blank will be used as a check on contamination of the samples.

The laboratory will be certified by the State of Illinois, and will adhere to a quality control program that meets specifications in 35 IAC 186, which governs the certification of laboratories.

Samples will be analyzed by standard methods according to USEPA SW-846 Test Methods for Evaluating Solid Waste. Quality control procedures listed in SW-846 for the methods that are used will be used by the laboratory.

## FEASIBILITY AND COST ANALYSIS

OF

## REMEDIAL ALTERNATIVES

MARCH 13, 2001

Dixon Marquette Cement Company
Dixon, Illinois

Prepared By Preston Engineering, Inc 4436 North Brady Avenue Davenport, Iowa 52806 Telephone: (319) 388-8288 Fax: (319) 388-9003

#### **EXECUTIVE SUMMARY**

Sixty four remediation technologies listed on the federal remediation technology roundtable were screened for applicability to Dixon Marquette Cement Company's landfill. Eleven technologies were considered theoretically feasible and were evaluated for installation at the landfill.

Installation of a RCRA Subtitle D Cap with Cap Enhancement was determined to be technically feasible. This type of technology would have a capital cost of \$2,000,000 to \$4,000,000.

No other technologies were feasible.

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#### I. INTRODUCTION

The Dixon Marquette Cement Plant (DMC) has disposed of waste cement kiln dust in an onsite landfill since 1984. The landfill appears to have been used as early as 1959 by previous plant owners. It is located in a portion of an old surface mine. The landfill is approximately 300 feet east of the Rock River at its closest point.

DMC desires to continue to use the landfill for disposal of waste cement kiln dust (CKD). However, because of inherent siting and design factors it is not technically or economically feasible to comply with all current landfill regulations. Therefore, it was necessary to evaluate alternatives that could be included in adjusted standards. Numerous sources were used to evaluate the costs and feasibility of various alternatives to the landfill regulations.

USEPA publication 540-R-00-002 "A Guide to Developing and Documenting Cost Estimates During the Feasibility Study" was used as a reference. The Federal Remediation Technology Roundtable website was used to screen various alternatives. A more detailed cost analysis was done for favorable alternatives.

The general alternatives considered were closing the landfill and hauling all new waste to the local landfill, hauling all new and previously disposed waste to the local landfill, and operating the landfill with adjusted standards. Options for collecting groundwater, treating groundwater, and hydraulic control were evaluated for the operating with adjusted standards alternative.

#### II. PRELIMINARY FEASIBILITY ANALYSIS

The Federal Remediation Technology Roundtable website (www.frtr.gov) contains a treatment technology screening matrix for many alternatives for treating waste, soil and groundwater. Table 3-2: Treatment Technologies Screening Matrix is included in the Appendix A. The inorganic contaminant column was selected because the CKD waste is inorganic and the leachate contaminants were primarily inorganic. The array of treatment technologies is rated by effectiveness. Technologies rated better for inorganic contaminants were selected for additional consideration. Background information was reviewed to determine if a technology could be used for the DMC site.

Many of the technologies rated better for inorganic technologies were determined to be ineffective, unproven, or expensive after review of background information. The following technologies were determined to be infeasible and were eliminated from further consideration. Technology descriptions are included in Appendix B.

- 4.6 Electrokinetic Separation
- 4.8 Soil Flushing
- 4.10 Solidification/Stabilization
- 4.17 Chemical Extraction
- 4.18 Chemical Reduction/Oxidation
- 4.20 Separation
- 4.21 Soil Washing
- 4.24 Solidification/Stabilization
- 4.40 Hydrofracturing
- 4.46 Passive/Reactive Treatment Walls
- 5.57 Deep Well Injection

The following technologies were determined to be theoretically feasible and were evaluated further to determine the site specific feasibility and cost at this facility. Technology descriptions are located in Appendix C.

- 4.30 Landfill Cap (Soil Containment Remediation)
- 4.31 Landfill Cap Enhancement (Soil Containment Remediation)
- 4.32 Excavation, Retrieval, and Off-Site Disposal (Soil Remediation)
- 4.36 Phytoremediation (In Situ Groundwater Remediation)
- 4.48 Constructed Wetlands (Ex Situ Groundwater Remediation)
- 4.49 Adsorption/Absorption (Ex Situ Groundwater Remediation)
- 4.52 Ion Exchange (Ex Situ Groundwater Remediation)
- 4.53 Precipitation/Coagulation/Flocculation (Ex Situ Groundwater Remediation)
- 4.40 Directional Wells (In Situ Groundwater Remediation)
- 4.58 Groundwater Pumping (Groundwater Containment)
- 4.59 Slurry Walls (In Situ Groundwater Remediation)

#### III. FEASIBILITY ANALYSIS OF SELECTED TECHNOLOGIES

Eleven technologies that were theoretically feasible were evaluated in more detail. The effectiveness at addressing all of the contaminants associated with the landfill as well as the cost was considered.

#### 4.30 Landfill Cap

The purpose of a cap is to prevent direct contact with waste, prevent wind blown transport of the waste, and to reduce the volume of precipitation that infiltrates through the waste and becomes leachate. The most common cap is a soil layer that has been compacted to former a barrier layer, which is then covered with a protective layer that supports vegetation. Two types of designs which are common are a RCRA Subtitle C Cap and a RCRA Subtitle D Cap.

A RCRA Subtitle D Cap is required for non-hazardous waste landfills. The Illinois regulation for Subtitle D landfills (811.314 Final Cover System) calls for a low permeability layer that has a minimum thickness of three feet and a permeability of less than 10<sup>-7</sup> cm/sec. A final protective cover is required that is three feet thick and is capable of supporting vegetation. Based on the technology summary, the rough industry cost is \$175k/acre acre for a RCRA Subtitle D. This works out to be a total cost of \$4,025,000 dollars for a 23-acre cap. However, because DMC has equipment to move soil it is anticipated the cost for a cap that meets Illinois requirements would cost less than \$2,000,000 if onsite soils can be used for the cap. A preliminary construction cost estimate for a landfill cap is included in Appendix D. The estimated cost for a landfill cap was \$1,852,000.

A RCRA Subtitle C Cap is a multilayered landfill cap that is suggested for use in RCRA hazardous waste applications. These caps have at least one synthetic liner in addition to the provisions of the Subtitle D cap. The synthetic liner will further reduce infiltration of water into the waste. The drawback is increased difficulty in construction and increased cost. The rough industry guideline cost is \$225k/acre. This would be a total of \$5,175,000 for a 23-acre cap. DMC does not have the ability to construct this type of cover and would need to use a landfill contractor. Therefore, the industry average is a reasonable estimate.

Construction of Subtitle D Cap installed in accordance with the IEPA rule 811.314 when the landfill is closed is a technologically feasible option.

#### 4.31 Landfill Cap Enhancements

The purpose of landfill cover enhancement is to reduce or eliminate percolation. Water harvesting increases runoff. Vegetation cover reduces soil moisture via plant uptake and evapotranspiration.

The anticipated design is likely to have at least a 20% slope which will result in a substantial amount of runoff. In addition, the landfill cap will need to be seeded with vegetation to control wind and water erosion of the soil cover. Landfill cap enhancements are anticipated as part of a normal soil cap construction and will not increase the cost of a Subtitle D cap.

Inclusion of landfill cap enhancements such as vegetation is a feasible and appropriate technology to apply to the anticipated Subtitle D cap.

#### 4.32 Excavation, Retrieval, and Off-Site Disposal

Contaminated material is removed and transported to permitted off-site treatment and disposal facilities. The technology guide provides a cost estimate of \$270 to \$460 per ton for disposal at a RCRA permitted facility.

The CKD landfill is estimated to contain 650,000 tons of waste. Disposal costs at the local permitted Subtitle D landfill are \$38/ton. The total cost-would be \$24,700,000 for disposal. There would be an additional cost of approximately \$300,000 for site restoration. The total cost for this technology is \$25,000,000.

The cost of offsite disposal is economically infeasible.

#### 4.36 Phytoremediation

Phytoremediation is a set of processes that uses plants to remove, transfer, stabilize and destroy organic/inorganic contamination in groundwater, surface water, and leachate. These mechanisms include enhanced rhizosphere biodegradation, hydraulic control, phyto-degradation, phyto-volatilization, phyto-accumulation, and phyto-stabilization.

The DMC site poses significant constraints on this technology. The primary contaminants are inorganic and thus are not amenable to biodegradation or volatilization. Much of the contamination is deep and in bedrock where roots are not likely to penetrate. Plants could pump shallow groundwater and accumulate contaminants in the shallow water near the Rock River. Planting trees along the river was investigated. However, there is very little area between the bluff overlooking the river and the river to plant trees. Currently, there is a lush stand of trees and vegetation along the bluff face that may very well attenuate contaminants in the shallow water table.

This technology was determined to be infeasible for this site because of site constraints.

#### 4.48 Constructed Wetlands

The constructed wetlands-based treatment technology uses natural geochemical and biological processes inherent in an artificial wetland ecosystem to accumulate and remove metals, explosives, and other contaminants from influent waters. The process can use a filtration or degradation process.

Although the technology incorporates principal components of wetland ecosystems, including organic soils, microbial fauna, algae, and vascular plants, microbial activity is responsible for most of the remediation. Metals are removed through ion exchange, adsorption, absorption, and precipitation with geochemical and microbial oxidation and reduction. Ion exchange occurs as metals in the water contact humic or other organic substances in the wetland. Wetlands constructed for this purpose often have little or no soil instead they have straw, manure or compost. Oxidation and reduction reactions catalyzed by bacteria that occur in the aerobic and anaerobic zones play a major role in precipitating metals as hydroxides and sulfides. Precipitated and adsorbed metals settle in quiescent ponds or are filtered out as water percolates through the medium or the plants.

The performance of wetland systems is not easily predicted nor is the cost. One estimate from the technical document is \$1.36 per thousand gallons. Up to three million gallons per day of contaminated groundwater is discharged. The cost for treating this volume of water would be \$4,080 per day or about \$1.5 million dollars per year.

This technology is not economically feasible and the technical feasibility is dubious.

#### 4.49 Adsorption/Absorption

In liquid absorption, solutes concentrate at the surface of the sorbent, thereby reducing their concentration in the bulk liquid phase. The most common adsorbent is granular activated carbon. Other adsorbents include activated alumina, forage sponge, lignin adsorption, sorption clays, and synthetic resins.

Adsorption does not work well for water soluble compounds and small molecules. Therefore, it would not be a good choice for inorganic contaminants such as dissolved solids. Cost estimates range from \$280/1000 gallons to \$340/1000 gallons. At three million gallons per day of contaminated groundwater, the cost could be over \$1,000,000 per day.

This technology will not work well for the dissolved constituents found in the groundwater at this site and the cost is prohibitory.

#### 4.52 Ion Exchange

Ion exchange removes ions from the aqueous phase by exchange with counter ions on the exchange medium. Ion exchange can remove dissolved metals from aqueous solutions. However, the ions removed are replaced with other ions with little change is dissolved solids content. Wastewater is also produced when the unit is regenerated. The cost is typically \$0.30 to \$0.80 per 1000 gallons treated. For three million gallons per day of contaminated groundwater, the cost would be \$900 to \$2,460 per day.

This technology would not significantly decrease the level of dissolved solids and the expense is significant.

#### 4.53 Precipitation/Coagulation/Flocculation

This process transforms dissolved contaminants into an insoluble solid, facilitating the contaminant's subsequent removal from the liquid phase by sedimentation or filtration. The process usually uses pH adjustment, addition of a chemical precipitant, and flocculation.

Precipitation of metals has long been the primary method of treating metal-laden industrial wastewaters. This involves the conversion of soluble heavy metal salts to insoluble salts that will precipitate. The precipitate can be removed from the treated water by physical methods such as ciarification (settling) and/or filtration. The precipitate or sludge must then be disposed in a landfill unless they have sufficient commercial value to allow recycling.

Cost for this process may be from \$0.30 to \$0.70 per 1,000 gallons plus \$0.50 per 1,000 gallons for sludge disposal. For a three million gallon per day system, the cost ranges from \$2,400 per day to \$3,750 per day.

This is a poor technology for treating groundwater under the DMC landfill. The groundwater has low heavy metal contamination. Adding chemicals will raise the dissolved solids content of the groundwater, which already has elevated levels. The annual cost is from \$876,000 to \$1,365,000 per year.

The technology is eliminated because of the high cost and questionable benefit.

#### 4.40 Directional Wells

Drilling techniques are used to position wells horizontally, or at an angle, to reach contaminants not accessible by direct vertical drilling.

Use of the technology may be limited by the potential for wells to collapse, requirement of specialized technology, difficulty in precisely locating wells, cost, and depths limited to 50 feet.

Cost estimates in the technology guide are \$20 to \$75 per foot. However, these costs appear to be for unconsolidated formations. Typically drilling in rock increase costs by a factor of 10. An analysis of leachate collection systems is included in Appendix E. The preliminary cost estimate for a horizontal well 2000 foot long is \$695,000 or \$350 per foot.

This technology appears feasible although there i some uncertainty of the percent of capture because of the variability of jointing in the bedrock.

#### 4.58 Cround water Pumping

Groundwater rumping is a commonly used groundwater remediation technology at contaminated sites. Possible objectives of groundwater pumping include removal of dissolved contaminants from the subsurface, and containment of contaminated groundwater to prevent migration.

Limitations of the technology include the length of time necessary to meet the remediation goal, the systems often fail to contain the contaminant plume, the cost of installation and operation is high, and biofouling of wells is common.

The cost for this type of system depends on the specific design and site conditions. A preliminary design is included in Appendix E. A system of 100 wells, 30 feet deep at DMC would cost over \$600,000 for construction. Operating costs were estimated to be \$100,000 per year.

This technology is feasible, although the actual capture of groundwater will depend on how effectively major joints are intersected. The logation of major joints is difficult to predict, making the percent of groundwater capture uncertain.

#### 4.59 Slurry Walls

These subsurface barriers consist of vertically excavated trenches filled with slurry. The slurry, usually a mixture of bentonite and water, hydraulically shores the trench to prevent collapse and retards ground water flow. Slurry walls are used to contain contaminated ground water, divert uncontaminated groundwater flow, and/or provide a barrier for the groundwater treatment system.

Slurry walls are typically placed at depths up to 100 feet and are generally 2 to 4 feet thick. The most effective application of a slurry wall for site remediation is to key the slurry wall 2 to 3 feet into a low permeability layer such as clay or bedrock. This keying-in provides for an effective foundation with minimal leakage potential.

The cost for a soil-bentonite wall in a soft to medium soil range is \$5 to \$7 per square foot. A recent bid for a local landfill resulted in a construction cost of approximately \$7 per square foot. These costs are based on excavation in soil. At the DMC site the slurry wall would have to be constructed in fractured dolomite and keyed into the Glenwood shale unit. Bedrock excavation is typically a factor of 10 higher than soil excavation. To circle the landfill with a 40-foot deep slurry wall would require a minimum of 160,000 square feet of slurry wall. At \$50 a square foot, the cost would be \$8,000,000.

The cost and difficulty in installing a slurry wall in 40 foot of rock makes this alternative infeasible.

## APPENDIX A

Treatment Technologies
Screening Matrix

Table 3-2: Treatment Technologies Screening Matrix

Rating Codes			Γ	<b></b> _	1		Ι	Г	Π	γ	<b></b>	T —		14	_	
■ - Better; O - Average;				ł										′		
△ - Worse. � - See Definition		ļ								l						
in Table 3-1a.		<b>=</b>							9/14	1.44	581		\$1	1940 Y	33.3	
Y - Yes; N - No.	}	reatment Train excludes off-gas treatment)	l	1			l				5					
F - Full; P - Pilot.		аţц	ı	ł			İ		S		Nonhalogenated SVOCs					
S - Solid; L - Liquid;	S II	Ę.	2			35			Š	S,	3	ည		基层		
V - Vapor.	evelopment Status	<u> </u>	Residuals Produced	l _		É			8	Halogenated VOCs	2	Halogenated SVOCs		100 H		
NA - Not Applicable	=	Train off-ga	Ę	O&M or Capital Intensive		iab Jity	2	يد	n3	5	2	'n		10	Radionuclides	
I - Inadequate.	Ē	s o	100	5	Availability	System Reliability Maintainability	Cleanup Time	Overall Cast	Nonhalogenated YOCs	ğ	g	384		norganics	Ē	Explosives
O&M - Operation &	g	Freatment (excludes	1 5	<u>ئ</u> ة	ap		5	=	- R	8	1	2	S	20	อแ	S
Maintenance; Cap - Capital; B -	1 %	\$ 5 E	, <u>š</u>	3 5	<b>1</b>	ste in	2	Š	5	<b>e</b>	5	9	Fuels	5	g	ē
Both	٥	150	<u> </u>	ÖĒ	₹	જે ¥ેં	<u> </u>	8	Z	Ξ	Z	=	ů.	5	α.	Ω
Soil, Sediment, and Slu	dge	e		عنود فدن ودر مشارکات	.,.,,											
				والمستندس	<u>ica</u>	<u>Trea</u>		ent								
4.1 Bioventing	F	N	N	N	�	•	1			<b>*</b>	�	�		Δ	*	I
4.2 Enhanced				[												
Bioremediation	F	N	N	0&м	�	•	•	4			•	4		•	•	
☐	1		l * `	Carr	•	•	*	*	_	_	*	*		<b>*</b>	4	_
☐ Anaerobic														•		
4.3 Land Treatment	F	N	N	N	<b>◆</b>	•	<b>*</b>	<b>*</b>				<b>*</b>		Δ	�	
4.4 Natural Attenuation	F	N	N	O&M	♦		$\triangle$	♦		•		•		<b>*</b>	4	
4.5 Phytoremediation		Ì														-
☐ ☐ Enhanced Rhizosphere			l													
Biodegradation	Р	N	N	N	Δ	0	Δ		0	0	0	0		Δ		
☐ ○ Phyto-accumulation					$\Delta$										Δ	
(? (?) Phyto-degradation	1				1										ĺ	
□ □ Phyto-stabilization			<u> </u>	L			<u> </u>		<u> </u>	<u> </u>				<u> </u>		
	In S	Situ P	hys	ical/C	her	nical '	<u> Fre</u>	atr	nen	<u>it</u>						
4.6 Electrokinetic	F	Y	L	0&M		0	0	0	0	0	0		_		^	_
Separation	1.	1	L	OWIVI	<b>452</b>	O			Γ			0	Δ		0	Δ
4.7 Fracturing						<del></del>										-
[] [] Blast-Enhanced	F	Y	NT.	0.34		_				_	٦	١, ١	اہرا	_		اہا
[] [] Lasagna Process	31	Y	N	O&M	$\triangle$	0	0	0	0	0	0	O	0	0	Δ	0
☐ ☐ Pneumatic Fracturing			L'	L												
4.8 Soil Flushing	Ŀ	1.4	Ţ,	00-14		^	,	_			_	_	~			<u> </u>
[11] Cosolvents Enhancement	1,	N	L	O&M	<b>3</b> 7	0		Δ			0	0	0		Δ	Δ
4.9 Soil Vapor Extraction	F	N	ſ.,	0&M		類	0		12		0	0		Δ	Δ	Δ
4,10							6									
Solidification/Stabilization	F	N	S	В	0	0	0	0								
(   [ ] In Situ Vitrification															_	
		3 In S	itn	There	nat	Treat	m.	n f	L	<b></b>	l,	<b></b> .	l	L	٠	L
	J.,	, ill ()	11,11	1 11011	1141	TICAL	ilic	ii Ç								

4.11 Thermally Enhanced Soil Vapor Extraction  □ Electrical Resistance Heating □ Radio Frequency/ Electromagnetic Heating □ Hot Air Injection  3.4 Ex Situ	P Ri	N	N	N	Δ	t (assu	Δ		0	0		•		Δ	Δ	Δ
4.12 Biopiles	F		T										-	<u> </u>	_	
	-	N	И	N			Ö		-		♦			*	$\stackrel{\triangle}{\cdot}$	
4.13 Composting	F	N	N	N			0				�	•		<b>*</b>	$\triangle$	
4.14 Fungal		* 1	·	000			١.	_								
<u>Biodegradation</u>	F	N	N	O&M	Δ	Δ			<b>*</b>	•	•	•	•	Δ	Δ	8
☐ ☐ White Rot Fungus							<u> </u>	ļ		<u> </u>		_				
4.15 Landfarming	F	N	N	N		<u> </u>					<b>*</b>	1			Δ	<b>*</b>
4.16 Slurry Phase	F	N	N	В	J	*0	0	0		_	•	•		<b>*</b>		
Biological Treatment	Υ.	14	14	Б					-		*	<b>*</b>		•		
3.5 Ex Situ Phy	sica	al/Cho	mi	cal Tr	eat	ment (	ass	un	iins	z ex	ca	vat	ion	)		
4.17 Chemical Extraction											Ī	Ī		_		
□□ Acid Extraction	F	Y	L	В	0	0		Δ	0	0			0		0	
□□ Solvent Extraction									į					_	_	
4.18 Chemical •		3.7				_					<u> </u>	_				
Reduction /Oxidation	F	Y	S	N		-		0	0	0	0	$ \circ $	0		Δ	$\triangle$
4.19 Dehalogenation										-	_	┢	-	-		
☐☐ Base-Catalyzed Decomposition	F	N	$ _{\mathbf{v}} $	В				Δ	0		0	_		_		
☐ ☐ Glycolate/Alkaline Polyethylene	r	IN	*	D	$\triangle$	Δ	Δ	<u> </u>		<u>ا</u>	10	0	Δ	$\triangle$	Δ	$\triangle$
Glycol (A/PEG)	<u> </u>	******				· <del></del>										
4.20 Separation																
☐☐ Gravity Separation	F	Y	S	0&M		0	120	0	0	0	0	lo			Δ	
☐ ☐ Magnetic Separation ☐ ☐ Sieving/Physical Separation		•	İ	_		_	l				] _	-		_		
CFC) Steving/Physical Separation		·	S					_								
4.21 Soil Washing	F	Y	L	В	100	0		0	0	0		-			Δ	23
4.22 Soil Vapor Extraction	F	N	니	N			_				_	<u> </u>	$\overline{}$			
The state of the s	P						0				<u></u>	0	$\circ$	$\triangle$	4	싘
4.23 Solar Detoxification	1	N	N	Cap.	0	0	0				=	#5		$\triangle$	$\triangle$	
4.24	ì											ł				
Solidification /Stabilization												•				
DD Bituminization											İ					
[] [] Emulsified Asphalt []] ] Modified Sulfur Cement	.,		ا ر													
[] [] Polyethylene Extrusion	F	N	S	Cap.				0	Δ	Δ	0	0	Δ			$\Delta$
Pozzolan/Portland Cement																
[] [] Radioactive Waste Solidification																
□ □ Sludge Stabilization												}				
[]] Soluble Phosphates																
[11] Vitrification/Molten Gluss	لـــا								ļ	L	<u> </u>	<u>L</u>				
3.6 Ex Situ	<u>T</u>	<u>ierma</u>	IT	<u> eatm</u>	ent	(assu	min	ge	xc	tva	tio	n)				
4.25 Hot Gas	Р	N	N	В					_		_	Ţ	_	_		
Decontamination	1	14	14	נו					47	4		^	Δ			

Mac v			_						<b></b> -		_			<del>,</del>		
4.26 Incineration			L		l											
☐ ☐ Circulating Bed Combustor ☐ ☐ Findized Bed	F	N	S	В		0		Δ					-			_
☐ ☐ Infrared Combustion	I.	14	V	D			-	4						44		
☐ Rotary Kiln			V													
4.27 Open Burn/Open	-		+		$\vdash$		├─	$\vdash$	-					-		
	F	N	S	В					Δ	Δ	Δ	Δ	Δ			
<u>Detonation</u>			<u> </u>									L				
4.28 Pyrolysis													-			ll
☐ ☐ Fluidized Bed	F	N	L	В	$\triangle$	I			0	0			0	Δ		ı
☐ ☐ Molten Salt Destruction	*	11	S	17	<b></b>	1			$\vee$		=		$  \cdot  $	2_3		1
□ □ Rotary Kiln							<u>L_</u>	L	L							
4.29 Thermal Desorption			L											*		
☐ ☐ High Temperature	F	Y	S	В		0		0		=				$\triangle$	$\triangle$	
☐ ☐ Low Temperature			o		1. 1											
	3.7 Containment															
4.30 Landfill Cap			T				Г									
□ □ Asphalt/Concrete Cap			L.,				١.		_	_		_		_		
□□ RCRA Subtitle C Cap	NA	N	NA	N					0	0	0	0	0	0		
□□ RCRA Subtitle D Cap																
4.31 Landfill Cap							<b>-</b>			-		_			-	-
Enhancements																
	NA	N	NA	N					0		0	0	0	0	$ \Delta $	
☐ ☐ Water Harvesting										1						
☐  Vegetative Cover			<u></u>				L	L		L	L					
3.8 Other Treatment																
4.32 Excavation, Retrieval,	NA	N	NA	N		=	10		0	0	0	0	0	0		
and Off-Site Disposal	LAZA	14	IVA	14		78	(4)			$  \cdot  $			$\Box$	$  \vee  $		0
Ground Water, Surfac	e W	ater.	ano	Leac	hat	te	*		·			L				
				Biolog		~~~~~	fme	ent								
4.33 Co-metabolic						2	Ť	1								
	P	N	N	O&M	Δ	•		1		<b>*</b>		<			Δ	ol
<u>Treatment</u>	<b> </b>						ļ		-		<u> </u>					
4.34 Enhanced																
Biodegradation																
□ □ Nitrate Enhancement	F	N	N	O&M		�	•	•		•	_	�			Λ.	
□ □ Oxygen Enhancement with Air	1	14	14	Occivi	-	**	*	W .		<b>*</b>	-	<b>₩</b>		$ \Delta $	$\triangle$	이
Sparging  Oxygen Enhancement with Hydrogen																
Peroxide							L									
4.35 Natural Attenuation	F	N	N	0&M		♦	•	•	<	•	4		*	Δ	Δ	Δ
4.36 Phytoremediation																
[11] Enhanced Rhizosphere	[						ļ	l								
Biodegradation	Р	N	N	N	0				0	0	0	0	0			اسرا
[1] Hydraulic Control		14	'	, ,		_	1	_	ľ	lٽ	$  \vee  $	\			Δ	
☐ ☐ Phyto-Degradation											'				l	
Phyto-Volatilization	لــا	~ .		<u> </u>	لــا		<u></u>	<u> </u>	L.	<b></b>	<u>L</u>	L	l	L		
				sical/C			Tre		mţ	ţ	*	·		-		]
4.37 Aeration	F	Y	V	N		0					0	0		Δ	Δ	Δ
4.38 Air Sparging	F	Y	V	N							$\triangle$	Δ		Δ	Δ	Δ
4.20 Disalumina	17	17	L	λī		$\sim$						-		Â		
4.39 Bioslurping	F	Y	l v	N		0	0		0	0				Δ	$ \Delta $	이
			4		L	·	<i>!</i>	I	I	ــــــــــــــــــــــــــــــــــــــ	J	I	l	J	li	lj

4.40 Directional Wells (enhancement)	F	N	NA	Сар.	Δ	0	趨	I	0	0	0	0	0	0	۵	0
4.41 Dual Phase Extraction	F	Y	L V	0&M		0	0	0		á	Δ	Δ		Δ	Δ	Δ
4.42 Fluid/Vapor Extraction	F	Y	L V	О&М		0	0	0			0	0	100	Δ	Δ	Δ
4.43 Hot Water or Steam Flushing/Stripping	P	Y	L V	Сар.		Δ		0	0	0	<b>G</b>		100	Δ	Δ	Δ
4.44 Hydrofracturing	P	Y	N	N	I			0	0	0	0	0	0	0	Δ	0
4 45 In-Well Air Stripping ☐ Circulating Wells	Р	Y	L V	Сар.			0	0		***	0	0		I	Δ	Δ
4.46 Passive /Rreactive Treatment Walls  Funnel and Gate	F	N	S	Сар.		I	Δ	I					0	<b>4</b>	Δ	
	3.1	1 Ex S	itu	Biolo	gica	ıl Tre	atn	en	t							
4.47 Biorcactors	F	N	S	Cap.		0	0				7	<b>*</b>		Δ	Δ	
4.48 Constructed Wetlands	F	N	S	Cap.	Δ	•	•	0	0	0	0	♦	0		Δ	
3.12 Ex Situ Pl	ıysi	cal/C	hen	ical T	rea	tmen	t (a	ssu	mi	ng	pui	no	ng	)		
4.49 Adsorption/ Absorption  □ Activated Alumina □ Forager Sponge □ Lignin Adsorption/ Sorptive Clays □ Synthetic Resins	Р	N	S	I	0	I	1	Δ	0	0	0	0	Δ		•	Δ
4.50 Air Stripping	F	N	L V	O&M			0				0	0	0	Δ	Δ	Δ
4.51 Granulated Activated Carbon (GAC)/Liquid Phase Carbon Adsorption	F	N	S	O&M	<b>D</b>			•					0	Δ	Δ	Δ
4.52 Ion Exchange	F	Y	S	Ν	7		0		Δ	Δ	Δ	$\triangle$	Δ		0	$\overline{\Delta}$
4.53 Precipitation/ Coagulation /Flocculation □ Coagulants and Flocculation	F	Y	S	N			0		Δ	Δ	Δ	Δ	Δ		0	I
4.54 Separation  □ □ Distillation □ □ Filtration/ Ultrafiltration /Microfiltration □ □ Freeze Crystallization □ □ Membrane Pervaporation □ □ Reverse Osmosis	F	Y	S	В				Δ		<b>748</b>			23	•	•	Δ
4.55 Sprinkler Irrigation	F	Y	S L	N	<b>34</b>	0	0							Δ	Δ	Δ
4.56 UV Oxidation □□ UV Photolysis	F	N	N	В			NA	0	14					Δ	Δ	
		3	.13	Conta	inr	nent										

4.57 Deep Well Injection	F	N	S L	N	12	0	NΑ		0	0	0	0	0	0	0	0
4.58 Ground Water Pumping □□ Surfactant Enhanced Recovery □□ Drawdown Pumping	F	N	L	В			NA	Δ	0	0	0	0	0	0	Δ	0
4.59 Slurry Walls	F	N	NA	Cap.					0	0	0	0	0	0	Δ	0
3.14 Air Emissions/Off-Gas Treatment																
4.60 Biofiltration	F	NA	S L	N	0	•	311	囊		<b>*</b>	•	•		Δ	NA	•
4.61 High Energy  Destruction  □ High Energy Corona  □ Tunable Hybrid Plasma Reactor	P	NA	N	I	Δ	Δ	ΝA	0	N				M	0	NΑ	Δ
4.62 Membrane Separation	P	NA	N	I	Δ	Δ	NΑ	0			0	0	0	Δ	NΑ	0
4.63 Oxidation  Catalytic Oxidation  In Internal Combustion Engine Oxidation  Thermal Oxidation  Ultraviolet Oxidation	F	NA	N	N	Ni.		NA	153	<b>1</b>	Æ	19		18	Δ	NΛ	0
4.64 Vapor Phase Carbon Adsorption  □ VOC Recovery and Recycle	F	NA	S	N		12	NA							0	NΑ	



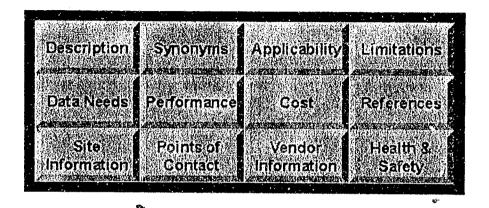
APPENDIX B

Infeasible Technologies

## 4.6 Electrokinetic Separation

(In Situ Soil Remediation Technology)

Previous Top Screen Table of Synonym Next
Section Page Matrix Contents List Section



Technology	Description
Soil, Sediment, and Sludge	
3,2 In Situ Physical/Chemical	Treatment
4.6 Electrokinetic Separation	The Electrokinetic Remediation (ER) process removes metals and organic contaminants from low permeability soil, mud, sludge, and marine dredging. ER uses electrochemical and electrokinetic processes to desorb, and then remove, metals and polar organics. This in situ soil processing technology is primarily a separation and removal technique for extracting contaminants from soils.

#### Description:



Figure 4-6: Typical In Situ Electrokinetic Separation System

The principle of electrokinetic remediation relies upon application of a low-intensity direct current through the soil between ceramic electrodes that are divided into a cathode array and an anode array. This mobilizes charged species, causing ions and water to move toward the electrodes. Metal ions, ammonium ions, and positively charged organic compounds move toward the cathode. Anions such as chloride, cyanide, fluoride, nitrate, and negatively charged organic compounds move toward the anode. The current creates an acid front at the anode and a base front at the cathode. This generation of acidic condition in situ may help to mobilize sorbed metal contaminants for transport to the collection system at the cathode.

The two primary mechanisms transport contaminants through the soil towards one or the other electrodes: electromigration and electroosmosis. In electromigration, charged particles are transported through the substrate. In contrast, electroosmosis is the movement of a liquid containing ions relative to a stationary charged surface. Of the two, electromigration is the main mechanism for the ER process. The direction and rate of movement of an ionic species will depend on its charge, both in magnitude and polarity, as well as the magnitude of the electroosmosis-induced flow velocity. Non-ionic species, both inorganic and organic, will also be transported along with the electroosmosis induced water flow.

Two approaches are taken during electrokinetic remediation: "Enhanced Removal" and "Treatment without Removal".

"Enhanced Removal" is achieved by electrokinetic transport of contaminants toward the polarized electrodes to concentrate the contaminants for subsequent removal and ex-situ treatment. Removal

of contaminants at the electrode may be accomplished by several means among which are: electroplating at the electrode; precipitation or co-precipitation at the electrode; pumping of water near the electrode; or complexing with ion exchange resins. Enhanced removal is widely used on remediation of soils contaminated metals.

"Treatment without Removal" is achieved by electro-osmotic transport of contaminants through treatment zones placed between electrodes. The polarity of the electrodes is reversed periodically, which reverses the direction of the contaminants back and forth through treatment zones. The frequency with which electrode polarity is reversed is determined by the rate of transport of contaminants through the soil. This approach can be used on in-situ remediation of soils contaminated with organic species.

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Synonyms:

Electrokinetics; Electromigration.

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Applicability:

Targeted contaminants for electrokinetics are heavy metals, anions, and polar organics in soil, mud, sledge, and marine dredging. Concentrations that can be treated range from a few parts per million (ppm) to tens of thousands ppm. Electrokinetics is most applicable in low permeability soils. Such soils are typically saturated and partially saturated clays and silt-clay mixtures, and are not readily drained.

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Limitations:

Factors that may limit the applicability and effectiveness of this process include:

- Effectiveness is sharply reduced for wastes with a moisture content of less than 10 percent.
   Maximum effectiveness occurs if the moisture content is between 14 and 18 percent.
- The presence of buried metallic or insulating material can induce variability in the electrical conductivity of the soil, therefore, the natural geologic spatial variability should be delineated. Additionally, deposits that exhibit very high electrical conductivity, such as ore deposits, cause the technique to be inefficient.
- Inert electrodes, such as carbon, graphite, or platinum, must be used so that no residue will
  be introduced into the treated soil mass. Metallic electrodes may dissolve as a result of
  electrolysis and introduce corrosive products into the soil mass.
- Electrokinetics is most effective in clays because of the negative surface charge of clay particles. However, the surface charge of the clay is altered by both charges in the pH of the pore fluid and the adsorption of contaminants. Extreme pH at the electrodes and reduction-oxidation changes induced by the process electrode reactions many inhibit ER's effectiveness, although acidic conditions (i.e., low pH) may help to remove metals.
- Oxidation/reduction reactions can form undesirable products (e.g., chlorine gas).

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Data Needs:

A detailed discussion of data elements is provided in <u>Subsection 2.2.1</u> (Data Requirements for Soil, Sediment and Sludge).

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Performance Data:

There have been few, if any, commercial applications of electrokinetic remediation in the United States. The electrokinetic technology has been operated for test and demonstration purposes at the pilot scale and at full scale at the following sites: (1) Louisiana State University, (2) Electrokinetics, Inc., (3) Geokinetics International, Inc., and (4) Battelle Memorial Institute. Geokinetics International, Inc.(GII) has successfully demonstrated the in situ electrokinetic remediation process in five field sites in Europe.

In 1996, a comprehensive demonstration study of lead extraction at a U.S.Army firing range in Louisiana was conducted by DoD's Small Business Innovative Research Program and

Electrokinetics, Inc. The EPA taking part in independent assessments of the results, found pilot-scale studies have demonstrated that concentrations of lead decreased to less than 300 mg/kg in 30 weeks of electrokinetic processing when the soils where originally contaminated as high as 4,500 mg/kg of lead.

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Cost:

Costs will vary with the amount of soil to be treated, the conductivity of the soil, the type of contaminant, the spacing of electrodes, and the type of process design employed. Ongoing pilot-scale studies using "real-world" soils indicate that the energy expenditures in extraction of metals from soils may be 500 kWh/m<sup>3</sup> or more at electrode spacing of 1.0m to 1.5m. Direct costs estimates of about \$15/m<sup>3</sup> for a suggested energy expenditure of \$0.03 per kilowatt hours, together with the cost of enhancement, could result in direct costs of \$50/m<sup>3</sup> or more. If no other efficient in situ technology is available to remediate fine-grained and heterogeneous subsurface deposits contaminated with metals, this technique would remain potentially competitive.

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References:

EPA, 1996. Recent Development for In Situ Treatment of Metal Contaminated Soil, EPA Office of Solid Waste & Emergency Response. Technology Innovative Office. Washington, DC.

USAEC, 1997. "In-situ Electrokinetic Remediation for Metal Contaminated Soils" in *Innovative Technology Demonstration, Evaluation and Transfer Activities, FY 96 Annual Report*, Report No. SFIM-AEC-ET-CR-97013, pp. 87-88.

U.S. DOE, 1995. "Electrokintic Remediation of Heavy Metals and Radionuclides," in *Technology Catalogue, Second Edition*, Office of Environmental Management Office of Technology Development, DOE/EM-0235, pp. 201-203.

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#### Site Information:

- Army Demo, U.S. Army Waterways Experiment Station, Baton Rouge Louisiana
- An Underground Storage Tank Spill
- DOE Demo, Oak Ridge K-25 facility, Tennessee
- DOE Demo: Sandia National Laboratories Chemical Waste Landfill
- EPA/DOE Demo: DOE Gaseous Diffusion Plant, Paducah, Kentucky
- DOE Demo, Old TNX Basin, Savannah River Site, South Carolina
- U.S. Army, firing range, LA
- Additional site information on the FRTR web site.

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**Points of Contact:** 

General FRTR Agency Contacts

**Technology Specific Web Sites:** 

Government Web Site is the Control of the Control o

### Non-Government Web Sites

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# Vendor Information:

A list of vendors offering In Situ Physical/Chemical Soil Treatment is available from the Vendor Information System for Innovative Treatment Technologies (VISITT) developed by U.S. Environmental Protection Agency (EPA).



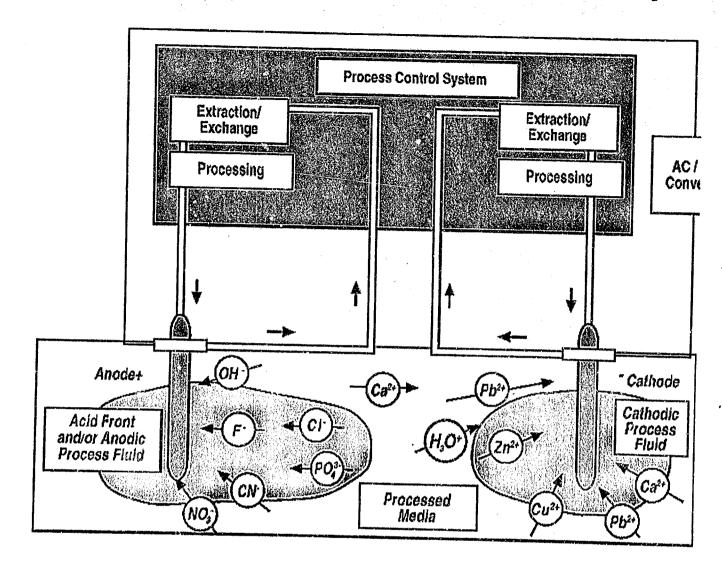
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To be added

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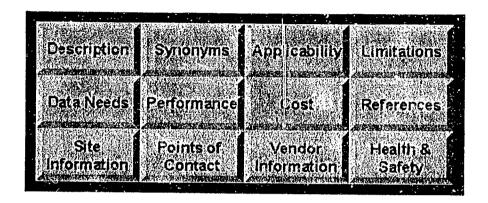
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# 4.8 Soil Flushing

In Site Soil Remediation Technology)

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Technology	Description
Soil, Sediment, and Sluc	ge
3.2 In Situ Physical/Che	mical Treatment
4.8 Soil Flushing	Water, or water containing an additive to enhance contaminant solubility, is applied to the soil or injected into the ground water to raise the water table into the contaminated soil zone. Contaminants are leached into the ground water, which is then extracted and treated.

Desc ption:



Figure 4-8: Typical Soil Flushing System

In situ soil flushing is the extraction of contaminants from the soil with water or other suitable aqueous solutions. Soil flushing is accomplished by passing the extraction fluid through in-place soils using an injection or infiltration process. Extraction fluids must be recovered from the underlying aquifer and, when possible, they are recycled.

## Cosolvent Enhancement

Cosolvent flushing involves injecting a solvent mixture (e.g., water plus a miscible organic solvent such as alcohol) into either vadose zone, saturated zone, or both to extract organic contaminants. Cosolvent flushing can be applied to soils to dissolve either the source of contamination or the contaminant plume emanating from it. The cosolvent mixture is normally injected upgradient of the contaminated area, and the solvent with dissolved contaminants is extracted downgradient and treated above ground.

Recovered ground water and flushing fluids with the desorbed contaminants may need treatment to meet appropriate discharge standards prior to recycle or release to local, publicly owned wastewater treatment works or receiving streams. To the maximum extent practical, recovered fluids should be reused in the flushing process. The separation of surfactants from recovered flushing fluid, for reuse in the process, is a major factor in the cost of soil flushing. Treatment of the recovered fluids results in process sludges and residual solids, such as spent carbon and spent ion exchange resin, which must be appropriately treated before disposal. Air emissions of volatile contaminants from recovered flushing fluids should be collected and treated, as appropriate, to meet applicable regulatory standards. Residual flushing additives in the soil may be a concern and should be

evaluated on a site-specific basis.

The duration of soil flushing process is generally short- to medium-term.

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Synonyms:

DSERTS Code: M12 (Soil Flushing).

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Applicability:

The target contaminant group for soil flushing is inorganics including radioactive contaminants. The technology can be used to treat VOCs, SVOCs, fuels, and pesticides, but it may be less cost-effective than alternative technologies for these contaminant groups. The addition of environmentally compatible surfactants may be used to increase the effective solubility of some organic compounds; however, the flushing solution may alter the physical/chemical properties of the soil system. The technology offers the potential for recovery of metals and can mobilize a wide range of organic and inorganic contaminants from coarse-grained soils.

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Limitations:

Factors that may limit the applicability and effectiveness of the process include:

- · Low permeability or heterogeneous soils are difficult to treat.
- Surfactants can adhere to soil and reduce effective soil porosity.
- Reactions of flushing fluids with soil can reduce contaminant mobility.
- The potential of washing the contaminant beyond the capture zone and the introduction of surfactants to the subsurface concern regulators. The technology should be used only where flushed contaminants and soil flushing fluid can be contained and recaptured.
- Aboveground separation and treatment costs for recovered fluids can drive the economics of the process.

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Data Needs:

A detailed discussion of these data elements is provided in <u>Subsection 2.2.1</u> (Data Requirements for Soil, Sediment, and Sludge). Treatability tests are required to determine the feasibility of the specific soil-flushing process being considered. Physical and chemical soil characterization parameters that should be established include soil permeability, soil structure, soil texture, soil porosity, moisture content, total organic carbon (TOC), cation exchange capacity (CEC), pH, and buffering capacity.

Contaminant characteristics that should be established include concentration, solubility, partition coefficient, solubility products, reduction potential, and complex stability constants. Soil and contaminant characteristics will determine the flushing fluids required, flushing fluid compatibility, and changes in flushing fluids with changes in contaminants.

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Performance Data:

Soil flushing is a developing technology that has had limited use in the United States. Typically, laboratory and field treatability studies must be performed under site-specific conditions before soil flushing is selected as the remedy of choice. To date, the technology has been selected as pirt of the source control remedy at 12 Superfund sites. This technology is currently operational at only one Superfund site; a second was scheduled to begin operation in 1991. EPA completed construction of a mobile soil-flushing system, the In Situ Contaminant/Treatment Unit, in 1988. This mobile soil-flushing system is designed for use at spills and uncontrolled hazardous waste sites. There has been very little commercial success with this technology.

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Cost:

The cost of soil flushing depends greatly on the type and concentration of surfactants used, if they are used at all. Rough estimates ranging from \$25 to \$250 per cubic yard have been reported.

Additional cost information can be found in the Hazardous, Toxic, and Radioactive Wastes (HTRW) Historical Cost Analysis System (HCAS) developed by Environmental Historical Cost Committee of Interagency Cost Estimation Group.



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References:

AATDF, 1997. Technology Practices Manual for Surfactants and Cosolvents, Technical Report, Document No. TR-97-2.

EPA, 1991. In Situ Soil Flushing, Engineering Bulletin, EPA/540/2-91/021.

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Sturges, S.G., Jr., P. McBeth, Jr., R.C. Pratt, 1992. "Performance of Soil Flushing and Groundwater Extraction at the United Chrome Superfund Site," *Journal of Hazardous Materials*, El Savior Science Pub., B.V., Amsterdam, Vol. 29, pp. 59-78.

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#### Site Information:

Laramic Tie Plant, WY

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**Points of Contact:** 



Technology Specific Web Sites:



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Vendor Information:

A list of vendors offering 'n Situ Physical/Chemical Soil Treatment is available from the Vendor Information System for Innovative Treatment Technologies (VISITT) developed by U.S. Environmental Protection Agency (EPA).



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Health and Safety:

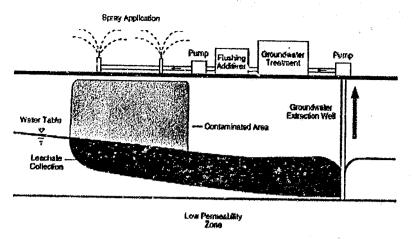


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4-5 OLP-3305 #22/04

# 4.10 Solidification/Stabilization

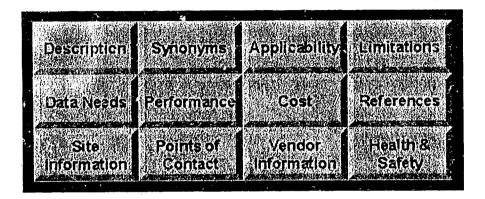
(In Situ Soil Remediation Technology)

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Technology	Description	
Soil, Sediment, and Sludge		
3.2 In Situ Physical/Chemical Treatment		
4.10 In Situ Solidification/ Stabilization	Contaminants are physically bound or enclosed within a stabilized mass (solidification), or chemical reactions are induced between the stabilizing agent and contaminants to reduce their mobility (stabilization).	

#### Description:

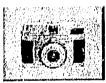


Figure 4-10a: Typical Auger/Caisson and Reagent/Injector Head Systems

Solidification/stabilization (S/S) reduces the mobility of hazardous substances and contaminants in the environment through both physical and chemical means. Unlike other remedial technologies, S/S seeks to trap or immobilize contaminants within their "host" medium (i.e., the soil, sand, and/or building materials that contain them), instead of removing them through chemical or physical treatment. Leachability testing is typically performed to measure the immobilization of contaminants. S/S techniques can be used alone or combined with other treatment and disposal methods to yield a product or material suitable for land disposal or, in other cases, that can be applied to beneficial use. These techniques have been used as both final and interim remedial measures.



Figure 4-10b: Typical In Situ Vitrification System

Auger/caisson systems and injector head systems are techniques used in soil S/S. They apply S/S agents to soils to trap or immobilize contaminants.

Bottom barriers are horizontal subsurface barriers that prevent vertical migration by providing a floor of impermeable material beneath the waste. The installation of a grout injection bottom barrier involves directional drilling with forced grout injection. Implementation of this technology is highly dependent on the physical properties of soil.

#### In Situ Vitrification (ISV)

In situ vitrification (ISV) is another in situ S/S process which uses an electric current to melt soil or other earthen materials at extremely high temperatures (1,600 to 2,000 °C or 2,900 to 3,650 °F) and thereby immobilize most inorganics and destroy organic pollutants by pyrolysis. Inorganic

pollutants are incorporated within the vitrified glass and crystalline mass. Water vapor and organic pyrolysis combustion products are captured in a hood, which draws the contaminants into an offges treatment system that removes particulates and other pollutants from the gas. The vitrification product is a chemically stable, leach-resistant, glass and crystalline material similar to obsidian or basalt rock. The process destroys and/or removes organic materials. Radionuclides and heavy metals are retained within the molten soil.

The timeframe for in situ S/S is short- to medium-term, while in situ ISV process is typically short-term.

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Synonyms:

**DSERTS** Codes:

M13 (Vitrification).

N11 (Solidification/Stabilization)

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Applicability:

The target contaminant group for in situ S/S is generally inorganics (including radionuclides).

The Auger/Caisson and Reagent/Injector Head Systems have limited effectiveness against SVOCs and pesticides and no expected effectiveness against VOCs; however, systems designed to be more effective in treating organics are being developed and tested.

The ISV process can destroy or remove organics and immobilize most inorganics in contaminated soils, sladge, or other earthen materials. The process has been tested on a broad range of VOCs and STOCs, other organics including dioxins and PCBs, and on most priority pollutant metals and radionuclides.

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Limitations:

Factors that may limit the applicability and effectiveness of the in situ S/S include:

- · Depth of contaminants may limit some types of application processes.
- Future usage of the site may "weather" the materials and affect ability to maintain immobilization of contaminants.
- Some processes result in a significant increase in volume (up to double the original volume).
- Certain wastes are incompatible with variations of this process. Treatability studies are generally required.
- Reagent delivery and effective mixing are more difficult than for ex situ applications.
- Like all in situ treatments, confirmatory sampling can be more difficult than for ex situ
  treatments.
- The solidified material may hinder future site use.
- Processing of contamination below the water table may require dewatering,

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Data Needs:

A detailed discussion of these data elements is provided in <u>Subsection 2.2.1</u> (Data Requirements for Soil, Sediment, and Sludge). Data needs include particle size, Atterberg limits, moisture content, metal concentrations, sulfate content, organic content, density, permeability, unconfined compressive strength, leachability, pH, and microstructure analysis. For ISV, a minimum alkali content in soil (sodium and potassium oxides) of 1.4 wt% is necessary to form glass. The composition of most soils is well within the range of processability.

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Performance Data:

Auger/Caisson and Reagent/Injector Head Systems processes are well demonstrated, can be applied to the most common site and waste types, require conventional materials handling equipment, and are available competitively from a number of vendors. Most reagents and additives are also widely available and relatively inexpensive industrial commodities.

Auger/Caisson and Reagent/Injector Head Systems processes have demonstrated the capability to reduce the mobility of contaminated waste by greater than 95%. The effects, over the long term, of weathering (e.g., freeze-thaw cycles, acid precipitation, and wind erosion), ground water infiltration, and physical disturbance associated with uncontrolled future land use can significantly affect the integrity of the stabilized mass and contaminant mobility in ways that cannot be predicted by laboratory tests.

There have been few, if any, commercial applications of ISV. The ISV process has been operated for test and demonstration purposes at the pilot scale and at full scale at the following sites: (1) Geosafe Corporation's test site, (2) DOE's Hanford Nuclear Reservation, (3) DOE's Oak Ridge National Laboratory, and (4) DOE's Idaho National Engineering Laboratory. More than 170 tests at various scales have been performed on a broad range of waste types in soils and sludge. A demonstration will take place at the Parsons/ETM site in Grand Ledge, Michigan, where the process is currently operating.

Process depths up to 6 meters (19 ft) have been achieved in relatively homogeneous soils. The achievable depth is limited under certain heterogeneous conditions.

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Cost:

Costs for Auger/Caisson and Reagent/Injector Head Systems processes vary widely according to materials or reagents used, their availability, project size, and chemical nature of contaminants (e.g., types and concentration levels for shallow applications). The in situ soil mixing/auger techniques average \$50 to \$80 per cubic meter (\$40 to \$60 per cubic yard) for the shallow applications and \$190 to \$330 per cubic meter (\$150 to \$250 per cubic yard) for the deeper applications.

The shallow soil mixing technique processes 36 to 72 metric tons (40 to 80 tons) per hour on average, and the deep soil mixing technique averages 18 to 45 metric tons (20 to 50 tons) per hour,

The major factor driving the selection process beyond basic waste compatibility is the availability of suitable reagents. Auger/Caisson and Reagent/Injector Head Systems processes require that potentially large volumes of bulk reagents and additives be transported to project sites. Transportation costs can dominate project economics and can quickly become uneconomical in cases where local or regional material sources are unavailable.

The cost for grout injection varies depending on site-specific conditions. Costs for drilling can range from \$50 to \$150/ft and grouting from \$50 to \$75/ft, not including mobilization, wash disposal, or adverse site condition expenses.

For ISV, average costs for treatability tests (all types) are \$25K plus analytical fees; for PCBs and dioxins, the cost is \$30K plus analytical. Equipment mobilization and demobilization costs are \$200K to \$300K combined. Vitrification operation cost varies with electricity costs, quantity of water, and depth of process.

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References:

California Base Closure Environmental Committee (CBCEC), 1994. Treatment Technologies Applications Matrix for Base Closure Activities, Revision 1, Technology Matching Process Action Team, November, 1994.

DOE, 1992. In Situ Vitrification, Technology Transfer Bulletin, prepared by Battelle's Pacific Northwest Laboratories for DOE, Richland, WA.

DOE, January 1992. "ISV Planning and Coordination," 1792 Technical Task Plan and Technical Task Description, TTP Reference No. RL-8568-PT.

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Technical Task Description, TTP Reference No. RL-8160-PT.

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EPA, 1989. SITE Program Demonstration Test International Waste Technologies In Situ Stabilization/Solidification Hialeah, Florida, Technology Evaluation Report, EPA RREL, Cincinnati, OH, EPA/540/5-89/004a.

EPA, 1989. Soliditech, Inc. Solidification, EPA RREL, series includes Technology Evaluation, Vol. I, EPA/540/5-89/005a; Technology Evaluation, Vol. II, EPA/540/5-89/005b, PB90-191768; Applications Analysis, EPA/540/A5-89/005; Technology Demonstration Summary, EPA/540/S5-89/005; and Demonstration Bulletin, EPA/540/M5-89/005.

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Federal Remediation Technologies Roundtable, 1995. Remediation Case Studies: Thermal Desorption, Soil Washing, and In Situ Vitrification, EPA/542/R-95/005.

 In Situ Vitrification at the Parsons Chemical/ETM Enterprises Superfund Site Grand Ledge, Michigan (interim results)

Federal Remediation Technologies Roundtable, 1997. Remediation Case Studies: Bioremediation and Vitrification, EPA/542/R-97/008.

- In Situ Vitrification at the Parsons Chemical/ETM Enterprises Superfund Site, Grand Ledge, Michigan
- In Situ Vitrification, U.S. Department of Energy, Hanford Site, Richland, Washington; Oak Ridge National Laboratory WAG 7; and Various Commercial Sites.

Rumer, R. and J.K., Mitchell (eds.), 1995. Assessment of Barrier Containment Technologies - a Comprehensive Treatment for Environmental Remediation Applications, International Containment Technology Workgroup, Baltimore.

Wiles, C.C., 1991. Treatment of Hazardous Waste with Solidification/Stabilization, EPA Report EPA/600/D-91/061.

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#### Site Information:

- EPA SITE Demo: Hialeah, FL
- EPA Demo: Geosafe Test Site, WA; Hanford Nuclear Reservation, WA
- · DOE Demo: Hanford Reservation, WA
- EPA Demo: Douglassville, PA
- Parsons Chemical/ETM Enterprises Superfund Site, Grand Ledge, MI (interim report)
- Parsons Chemical/ETM Enterprises Superfund Site, Grand Ledge, MI (final report)
- Hanford Site, Richland, WA; Oak Ridge National Lab WAG 7
- Additional site information on the FRTR web site

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Vendor Information:

A list of vendors offering In Situ Physical/Chemical Soil Treatment is available from the Vendor Information System for Innovative Treatment Technologies (VISITT) developed by U.S. Environmental Protection Agency (EPA).



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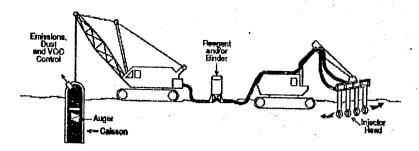
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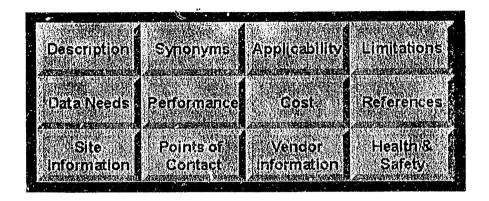


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# 4.17 Chemical Extraction

(Ex Situ Soil Remediation Technology)

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Technology	Description	
Soil, Sediment, and Sludge		
3.5 Ex Situ Physical/Chemical Treatment (assuming excavation)		
4.17 Chemical Extraction	Waste contaminated soil and extractant are mixed in an extractor, dissolving the contaminants.  The extracted solution is then placed in a separator, where the contaminants and extractant are separated for treatment and further use.	

#### Description:



Figure 4-17:
Typical Solvent Extraction
Process

Chemical extraction does not destroy wastes but is a means of separating hazardous contaminants from soils, sludges, and rediments, thereby reducing the volume of the hazardous waste that must be treated. The technology uses an Extracting chemical and differs from soil washing, which generally uses water or water with wash-improving additives. Commercial-scale units are in operation. They vary in regard to the Chemical employed, type of equipment used, and mode of operation.

Physical separation steps are often used before chemical extraction to grade the soil into coarse and fine fractions, with the assumption that the fines contain most of the contamination. Physical separation can also enhance the kinetics of extraction by separating out particulate heavy metals, if these are present in the soil.

#### > Acid Extraction

Acid can also be used as the extractant. Acid extraction uses hydrochloric acid to extract heavy metal contaminants from soils. In this process, soils are first screened to remove coarse solids. Hydrochloric acid is then introduced into the soil in the extraction unit. The residence time in the unit varies depending on the soil type, contaminants, and contaminant concentrations, but generally ranges between 10 and 40 minutes. The soil-extractant mixture is continuously pumped out of the mixing tank, and the soil and extractant are separated using hydrocyclones.

When extraction is complete, the solids are transferred to the rinse system. The soils are rinsed with water to remove entrained acid and metals. The extraction solution and rinse waters are regenerated

using comercially available precipitants, such as sodium hydroxide, lime, or other proprietary formulations, along with a flocculent that removes the metals and reforms the acid. The heavy metals are concentrated in a form potentially suitable for recovery. During the final step, the soils are dewatered and mixed with lime and fertilizer to neutralize any residual acid.

#### Solvent Extraction

Solvent extraction is a common form of chemical extraction using organic solvent as the extractant. It is commonly used in combination with other technologies, such as solidification/stabilization, incineration, or soil washing, depending upon site-specific conditions. Solvent extraction also can be used as a stand alone technology in some instances. Organically bound metals can be extracted along with the target organic contaminants, thereby creating residuals with special handling requirements. Traces of solvent may remain within the treated soil matrix, so the toxicity of the solvent is an important consideration. The treated media are usually returned to the site after having met Best Demonstrated Available Technology (BDAT) and other standards.

The duration of operations and maintenance for chemical extraction is medium-term.

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Synonyms:

**DSERTS Codes:** 

N16 (Acid Extraction)
N17 (Solvent Extraction)

Backto Ton

Applicability:

Solvent extraction has been shown to be effective in treating sediments, sludges, and soils containing primarily organic contaminants such as PCBs, VOCs, halogenated solvents, and petroleum wastes. The process has been shown to be applicable for the separation of the organic contaminants in paint wastes, synthetic rubber process wastes, coal tar wastes, drilling muds, wood-treating wastes, separation sludges, pesticide/insecticide wastes, and petroleum refinery oily wastes.

Acid extraction is suitable to treat sediments, sludges, and soils contaminated by heavy metals.

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Lire itations:

Factors that may limit the applicability and effectiveness of the process include:

- Some soil types and moisture content levels will adversely impact process performance.
- Higher clay content may reduce extraction efficiency and require longer contact times.
- Organically bound metals can be extracted along with the target organic pollutants, which
  restricts handling of the residuals.
- The presence of detergents and emulsifiers can unfavorably influence the extraction performance.
- Traces of solvent may remain in the treated solids; the toxicity of the solvent is an important consideration.
- Solvent extraction is generally least effective on very high molecular weight organic and very hydrophilic substances.
- After acid extraction, any residual acid in treated soil needs to be neutralized.
- Capital costs can be relatively high and the technology may be more economical at larger sites.
- Meeting highly stringent heavy metals criteria (e.g., passing the California WET test) may
  prove uneconomical.

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Data Needs:

A detailed discussion of these data elements is provided in <u>Subsection 2.2.1</u> (Data Requirements for Soil, Sediment, and Sludge). It is important to determine whether mass transfer or equilibrium will be controlling. The controlling factor is critical to the design of the unit and to the determination of whether the technology is appropriate for the waste.

Soil properties that should be determined include particle size; pH; partition coefficient; cation exchange capacity; organic content; TCLP; moisture content; and the presence of metals, volatiles, clays, and complex waste mixtures.

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Performance Data:

The performance data currently available are mostly from Resource Conservation Company (RCC). The ability of RCC's full-scale B.E.S.T. The process to separate oily feedstock into product fractions was evaluated by EPA at the General Refining Superfund site near Savannah, Georgia, in February 1987. The treated soils from this unit were backfilled to the site, product oil was recycled as a fuel oil blend, and the recovered water was pH-adjusted and transported to a local industrial wastewater treatment facility.

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Cost:

Cost estimates for this technology range from \$110 to \$440 per metric ton (\$100 to \$400 per ton), depending on the volume of soil treated.

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References:

California Base Closure Environmental Committee (CBCEC), 1994. *Treatment Technologies Applications Matrix for Base Closure Activities, Revision 1*, Technology Matching Process Action Team, November, 1994.

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Federal Remediation Technologies Roundtable, 1998. Remediation Case Studies: Ex Situ Soil Treatment Technologies (Bioremediation, Solvent Extraction, Thermal Desorption), EPA/542/R-98/011.

Solvent Extraction at the Sparrevoln Long Range Radar Station, Alaska

Raghavan, R., D.H. Dietz, and E. Coles, 1988. Cleaning Excavated Soil Using Extraction Agents: A State-of-the-Art Review, EPA Releases Control Branch, Edison, NJ, EPA Report EPA 600/2-89/034.

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#### Site Information:

- EPA Removal Action: Traband Warehouse PCBs, OK
- EPA Removal Action: Stanford Pesticide Site No. 1, AZ
- EPA Remedial Action: Palmetto Wood Preserving, SC
- EPA Removal Action: PBM Enterprises, MI
- EPA Removal Action: Zhiegner Refining Company
- EPA Demo: Midwest, California, Australia
- EPA Demo: Grand Calumet River Site, IL
- EPA SITE Demo: Research Facility, Edison, NJ
- United Creosoting, Conroe, TX
- DOE pilot-scale test, Idaho National Engineering Laboratory
- · SUPERFUND Remedial Action, Arrowhead Refinery Co.
- EPA SITE Demo: Naval Air Station North Island Site 4, San Diego, CA
- EPA Removal Action: General Refining Company, GA
- EPA Removal Action: Vineland Chemical Company, NJ
- EPA Removal Action: Avtex Fibers, VA
- EPA & Navy Demo: EPA Lab, NJ
- EPA Demo; Douglassville, PA
- EPA Demo
- Star Enterprise, Port Arthur, TX
- EPA Demo: New Bedford Harbor, MA & O'Connor Site, ME
- Sparrevohn Long Range Radar Station, AK
- Additional site information on the FRTR web site

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**Points of Contact:** 

General FRTR Agency Contacts

**Technology Specific Web Sites:** 

Non-Government Web Sites

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#### Vendor Information:

A list of vendors offering Ex Situ Physical/Chemical Soil Treatment is available from the Vendor Information System for Innovative Treatment Technologies (VISITT) developed by U.S. Environmental Protection Agency (EPA).



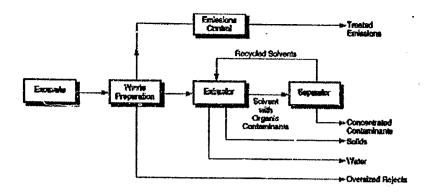
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# 4.18 Chemical Reduction/Oxidation

(Ex Situ Soil Remediation Technology)

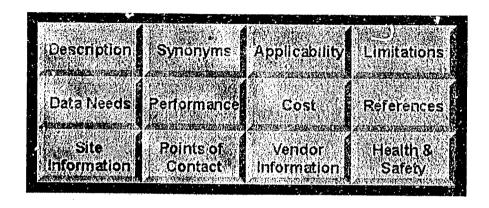
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Technology	Description
Soil, Sediment, and Sludge	
3.5 Ex Situ Physical/Chemi	ical Treatment (assuming excavation)
4.18 Chemical Reduction/Oxidation	Reduction/oxidation chemically converts hazardous contaminants to non-hazardous or less toxic compounds that are more stable, less mobile, and/or inert. The oxidizing agents most commonly used are ozone, hydrogen peroxide, hypoclylorites, chlorine, and chlorine dioxide.

#### Description:

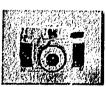


Figure 4-18:
Typical Chemical
Reduction/Oxidation Process

Reduction/oxidation (Redox) reactions chemically convert hazardous contaminants to nonhazardous or less toxic compounds that are more stable, less mobile, and/or inert. Redox reactions involve the transfer of electrons from one compound to another. Specifically, one reactant is oxidized (loses electrons) and one is reduced (gains electrons). The oxidizing agents most commonly used for treatment of hazardous contaminants are ozone, hydrogen peroxide, hypochlorites, chlorine, and chlorine dioxide. Chemical redox is a full-scale, well-established technology used for disinfection of drinking water and wastewater, and it is a common treatment for cyanide wastes. Enhanced systems are now being used more frequently to treat contaminants in soils.

Chemical reduction/oxidation is a short- to medium-term technology.

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DSERTS Code: N13 (Chemical Reduction/Oxidation).

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The target contaminant group for chemical redox is inorganics. The technology can be used but may be less effective against nonhalogenated VOCs and SVOCs, fuel hydrocarbons, and pesticides.

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Limitations:

Synonyms:

Applicability:

Factors that may limit the applicability and effectiveness of the process include:

- Incomplete oxidation or formation of intermediate contaminants may occur depending upon the contaminants and oxidizing agents used.
- The process is not cost-effective for high contaminant concentrations because of the large amounts of oxidizing agent required.
- Oil and grease in the media should be minimized to optimize process efficiency.

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Data Needs:

A detailed discussion of these data elements is provided in <u>Subsection 2.2.1</u> (Data Requirements for Soil, Sediment, and Sludge). Treatability tests should be conducted to identify parameters such as water, alkaline metals, and humus content in the soils; the presence of multiple phases; and total organic halides that could affect processing time and cost.

Performance Data:

Chemical redox is a full-scale, well-established technology used for disinfection of drinking water and wastewater, and it is a common treatment for cyanide (oxidation) and chromium (reduction of Cr (VI) to Cr (III) prior to precipitation) wastes. Enhanced systems are now being used more frequently to treat hazardous wastes in soils.

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Cost:

Estimated costs range from \$190 to \$660 per cubic meter (\$150 to \$500 per cubic yard).

Additional cost information can be found in the Hazardous, Toxic, and Radioactive Wastes (HTRW) Historical Cost Analysis System (HCAS) developed by Environmental Historical Cost Committee of Interagency Cost Estimation Group.



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References:

California Base Closure Environmental Committee (CBCEC), 1994. *Treatment Technologies Applications Matrix for Base Closure Activities, Revision 1*, Technology Matching Process Action Team, November, 1994.

EPA, 1993. Lawrence Livermore National Laboratory Superfund Site, Project Summary, EPA/540/SR-93/516.

EPA, 1991. *Chemical Oxidation Treatment*, Engineering Bulletin, EPA, OERR and ORD, Washington, DC, EPA/530/2-91/025.

Mayer, G., W. Bellamy, N. Ziemba, and L.A. Otis, 15-17 May 1990. "Conceptual Cost Evaluation of Volatile Organic Compound Treatment by Advanced Ozone Oxidation," Second Forum on Imposative Hazardous Waste Treatment Technologies: Domestic and International, Philadelphia, PA, EPA, Washington, DC, EPA Report EPA/2-90/010.

USAEC, 1997. "Catalyzed Hydrogen Peroxide Treatment of 2, 4, 6-Trinitrotoluene in Soils" in Innovative Technology Demonstration, Evaluation and Transfer Activities, FY 96 Annual Report, Report No. SFIM-AEC-ET-CR-97013, pp. 77-78.

USAEC, 1997. "Remediation of Chemical Agent Contaminated Soils Using Peroxysulfate" in *Innovative Technology Demonstration, Evaluation and Transfer Activities, FY 96 Annual Report*, Report No. SFIM-AEC-ET-CR-97013, pp. 93-94.

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Site Information:

- EPA Demo: Coleman-Evans Site, FL
- Additional site information on the FRTR web site

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**Points of Contact:** 

General FRTR Agency Contacts

**Technology Specific Web Site:** 



Non-Government Web Sites

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#### Vendor Information:

A list of vendors offering Ex Situ Physical/Chemical Soil Treatment is available from the Vendor Information System for Innovative Treatment Technologies (VISITT) developed by U.S. Environmental Protection Agency (EPA).



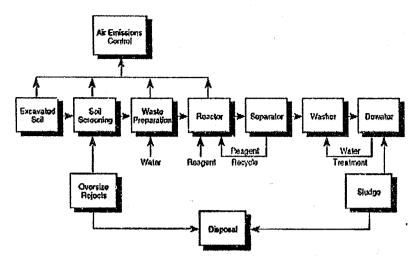
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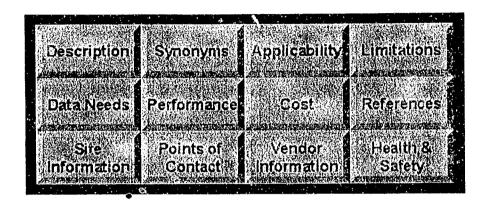
# 4.20 Separation

(Ex Situ Soil Remediation Technology)

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Technology	Description
Soil, Sediment, and Sludge	
3.5 Ex Situ Physical/Chemical Treatment	
4.20 Separation	Separation techniques concentrate contaminated solids through physical and chemical means. These processes seek to detach contaminants from their medium (i.e., the soil, sand, and/or binding material that contains them).

#### Description:

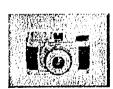


Figure 4-20: Typical Gravity **Separation** System

Ex situ separation can be performed by many processes. Gravity separation and sieving/physical separation are two well-developed processes that have long been primary methods for treating municipal wastewaters. Magnetic separation, on the other hand, is a much newer separation process that is still being tested.

#### > Gravity Separation

Gravity separation is a solid/liquid separation process, which relies on a density difference between the phases. Equipment size and effectiveness of gravity separation depends on the solids settling velocity, which is a function of the particles size, density difference, fluid viscosity, and particle concentration (hindered settling). Gravity separation is also used for removing immiscible oil phases, and for classification where particles of different sizes are separated. It is often preceded by coagulation and flocculation to increase particle size, thereby allowing removal of fine particles.

#### > Magnetic Separation

Magnetic separation is used to extract slightly magnetic radioactive particles from host materials such as water, soil, or air. All uranium and plutonium compounds are slightly magnetic while most host materials are nonmagnetic. The process operates by passing contaminated fluid or slurry through a magnetized volume. The magnetized volume contains a magnetic matrix material such as steel wool that extracts the slightly magnetic contamination particles from the slurry.

Sieving/Physical Separation

Sieving and physical separation processes use different size sieves and screens to effectively concentrate contaminants into smaller volumes. Physical separation is based on the fact that most organic and inorganic contaminants tend to bind, either chemically or physically, to the fine (i.e., clay and silt) fraction of a soil. The clay and silt soil particles are, in turn, physically bound to the coarser sand and gravel particles by compaction and adhesion. Thus, separating the fine clay and silt particles from the coarser sand and gravel soil particles would effectively concentrate the contaminants into a smaller volume of soil that could then be further treated or disposed.

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Synonyms:

NA

Applicability:

The target contaminant groups for ex situ separation processes are SVOCs, fuels, and inorganics (including radionuclides). The technologies can be used on selected VOCs and pesticides. Magnetic separation is specifically used on heavy metals, radionuclides, and magnetic radioactive particles, such as uranium and plutonium compounds.

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Physical separation often precedes chemical extraction treatment based on the assumption that most of the contamination is tied to the finer soil fraction, which alone may need to be treated. Separation is also useful when heavy metal contaminants occur as particulates (e.g., in small-arms ranges). One advantage of physical separation processes is that high throughputs can be achieved with relatively small equipment.

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Limitations:

Factors that may limit the applicability and effectiveness of these processes include:

- High clay and moisture content will increase treatment cost.
- Gravity separation processes rely on a difference in the solids and liquid phase densities.
   Specific gravity of particles will effect settling rate and process efficiency. Additionally, settling velocity is dependent on the viscosity of the suspending fluid, which must be known to estimate process efficiency and to size equipment.
- Special measures may be required to mitigate odor problems, resulting from organic sludge that undergoes septic conditions.

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Data Needs:

A detailed discussion of data elements is provided in <u>Subsection 2.2.1</u> (Data Requirements for Soil, Sediment, and Sludge). In addition, particle size distribution; soil type, physical form, handling properties, and moisture content; contaminant type and concentration; texture; and organic content need to be investigated.

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Performance Data:

Gravity separation and sieving/physical separation are full-scale, well-established technologies used mostly for treatment of wastewater and contaminated soil, sediment, and sludge. Magnetic separation is a promising new technique used to remove radioactive contaminants from soils. It has recently been tested at the bench-scale level at DOE sites.

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Cost:

NA

Additional cost information can be found in the Hazardous, Toxic, and Radioactive Wastes (HTRW) Historical Cost Analysis System (HCAS) developed by Environmental Historical Cost Committee of Interagency Cost Estimation Group.

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References:

Battelle Memorial Institute, 1995. ReOpt. V3.1, by Battelle Memorial Institute for DOE under

Contract DE/AC06/76RLO 1830.

Coyne Prenger, 1995. High Gradient Magnetic Separation (HGMS), Los Alamos National

Laboratory (LANL), LALP-94-264.

DOE, April 1995. Technology Catalogue, Second Edition, Office of Environmental Management

& Office of Technology Development, DOE/EM-0235.

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#### Site Information:

- Los Alamos National Laboratory (LANL)
- DOE Demo: INEL, ID
- EPA Demo: Iron Mountain Mine Site, CA
- Twin Cities AAP, New Brighton, MN
- Additional site information on the FRTR web site

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Points of Contact:



**Technology Specific Web Sites:** 



#### Vendor Information:

A list of vendors offering Ex Situ Physical/Chemical Soil Treatment is available from the Vendor Information System for Innovative Treatment Technologies (VISITT) developed by U.S. Environmental Protection Agency (EPA).



Health and Safety:



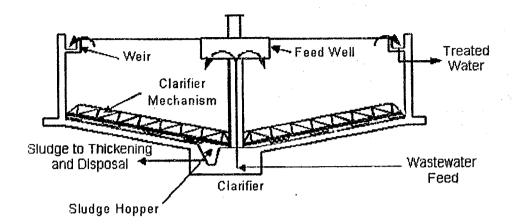
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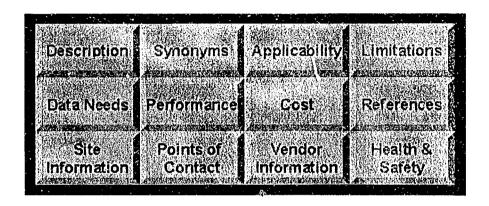
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# 4.21 Soil Washing

(Ex Situ Soil Remediation Technology)

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Technology	Description
Soil, Sediment, and Slud	ge
3.5 Ex Situ Physical/Che	mical Treatment (assuming excavation)
4.21 Soil Washing	Contaminants sorbed onto fine soil particles are separated from bulk soil in an aqueous-based system on the basis of particle size. The wash water may be augmented with a basic leaching agent, surfactant, pH adjustment, or chelating agent to help remove organics and heavy metals.

Description:



Figure 4-21: Typical Soil Washing Process

Soil washing is a water-based process for scrubbing soils ex situ to remove contaminants. The process removes contaminants from soils in one of two ways:

- By dissolving or suspending them in the wash solution (which can be sustained by chemical manipulation of pH for a period of time).
- By concentrating them into a smaller volume of soil through particle size separation, gravity separation, and attrition scrubbing (similar to those techniques used in sand and gravel operations).

Soil washing systems incorporating most of the removal techniques offer the greatest promise for application to soils contaminated with a wide variety of heavy metal, radionuclides, and organic contaminants. Commercialization of the process, however, is not yet extensive.

The concept of reducing soil contamination through the use of particle size separation is based on the finding that most organic and inorganic contaminants tend to bind, either chemically or physically, to clay, silt, and organic soil particles. The silt and clay, in turn, are attached to sand and gravel particles by physical processes, primarily compaction and adhesion. Washing processes that separate the fine (small) clay and silt particles from the coarser sand and gravel soil particles effectively separate and concentrate the contaminants into a smaller volume of soil that can be further treated or disposed of. Gravity separation is effective for removing high or low specific gravity particles such as heavy metal-containing compounds (lead, radium oxide, etc.). Attrition scrubbing removes adherent contaminant films from coarser particles. However, attrition washing can increase the fines in soils processed. The clean, larger fraction can be returned to the site for continued use.

Complex mixture of contaminants in the soil (such as a mixture of metals, nonvolatile organics, and SVOCs) and heterogeneous contaminant compositions throughout the soil mixture make it difficult to formulate a single suitable washing solution that will consistently and reliably remove all of the different types of contaminants. for these cases, sequential washing, using different wash formulations and/or different soil to wash fluid ratios, may be required.

Soil washing is generally considered a media transfer technology. The contaminated water generated from soil washing are treated with the technology(s) suitable for the contaminants.

The duration of soil washing is typically short- to medium-term.

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Synonyms:

**DSERTS Code:** N15 (Soil Washing).

Applicability:

The target contaminant groups for soil washing are SVOCs, fuels, and heavy metals. The technology can be used on selected VOCs and pesticides. The technology offers the ability for recovery of metals and can clean a wide range of organic and inorganic contaminants from coarse-

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grained soils.

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Limitations:

Factors that may limit the applicability and effectiveness of the process include:

- Complex waste mixtures (e.g., metals with organics) make formulating washing fluid difficult.
- High humic content in soil may require pretreatment.
- The aqueous stream will require treatment at demobilization.
- Additional treatment steps may be required to address hazardous levels of washing solvent remaining in the treated residuals.
- It may be difficult to remove organics adsorbed onto clay-size particles.

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Data Needs:

A detailed discussion of these data elements is provided in <u>Subsection 2.2.1</u> (Data Requirements for Soil, Sediment, and Sludge). Particle size distribution (6.24 to 2 mm optimum range); soil type, physical form, handling properties, and moisture content; contaminant type and concentration; texture; organic content; cation exchange capacity; pH and buffering capacity. A complete bench scale treatability study should always be completed before applying this technology as a remedial solution.

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Performance Data:

At the present time, soil washing is used extensively in Europe but has had limited use in the United States. During 1986-1989, the technology was one of the selected source control remedies at eight Superfund sites.

Soil washing provides a cost effective and environmentally proactive alternative to stabilization and landfilling. Two pilot scale demonstrations were carried out at Fort Polk, Louisiana in 1996. These employed commercially available unit processes - physical separation/acid leaching systems. The system employed acetic acid as the leaching agent, and the other, hydrochloric acid. Input soil had a lead content of approximately 3500 mg/kg. The hydrochloric acid system was most effective. Processed soil had total lead concentration of 200 mg/kg and TCLP levels for lead of approximately 2 mg/L. The through put rate was approximately 6 tons per hour. Choice of acid leaching agent is a function of specific soil chemistry and degree of solubility required.

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Cost:

The average cost for use of this technology, including excavation, is approximately \$170 per ton, depending on site specific conditions and the target waste quantity and concentration.

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References:

Battelle, 1997. Physical Separation and Acid Leaching: A Demonstration of Small-Arms Range Remediation at Fort Polk, Louisiana. Final report prepared for Naval Facilities Engineering Service Center (NFESC) and U.S. Army Environmental Center under contract with NFESC, Port Hueneme, CA.

Battelle, 1997. Physical Separation and Acid Leaching: A Demonstration of Small-Arms Range Remediation at Fort Polk, Louisiana. Implementation Guidance Handbook. Prepared for Naval Facilities Engineering Service Center (NFESC) and U.S. Army Environmental Center under contract with NFESC, Port Hueneme, CA.

BDM, 1997. Demonstration of Physical Separation/Leaching Methods for the Remediation of Heavy Metals-Contaminated Soils at Small-Arms Range Final Report. Prepared for U.S. Army Environmental Center (USAEC) and Naval Facilities Engineering Service Center under contract with USAEC, Aberdeen Proving Ground, MD.

BDM, 1997. Demonstration of Physical Separation/Leaching Methods for the Remediation of Heavy Metals-Contaminated Soils at Small-Arms Range Worldwide Search Report. Prepared for U.S. Army Environmental Center (USAEC) and Naval Facilities Engineering Service Center under contract with USAEC, Aberdeen Froving Ground, MD.

California Base Closure Environmental Committee (CBCEC), 1994. Treatment Technologies Applications Matrix for Base Closure Activities, Revision 1, Technology Matching Process Action Team, November, 1994.

EPA, 1989. Innovative Technology: Soil Washing, OSWER Directive 9200.5-250FS.

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EPA, 1990. Soil Washing Treatment, Engineering Bulletin, EPA, OERR, Washington, DC, EPA/540/2-90/017. Available from NTIS, Springfield, VA, Order No. PB91-228056.

EPA, 1991. *Blotrol Soll Washing System*, EPA RREL, series includes Technology Evaluation Vol. I, EPA/540/5-91/003a, PB92-115310; Technology Evaluation Vol. II, Part A, EPA/540/5-91/003b, PB92-115328; Technology Evaluation Vol. II, Part B, EPA/540/5-91/003c, PB92-115336; Applications Analysis, EPA/540/A5-91/003; Technology Demonstration Summary, EPA/540/S5-91/003; and Demonstration Bulletin, EPA/540/M5-91/003.

EPA, 1992. A Cuizen's Guide to Soil Washing, EPA, OSWER, Washington, DC, EPA/542/F-92/003.

EPA, 1992. Bergmann USA Soll/Sedlment Washing System, EPA RREL, Demonstration Bulletin, EPA/540/MR-92/075.

EPA, 1993. Bescorp Soil Washing System Battery Printerprises Site Brice Environmental Services, Inc., EPA RREL, Demonstration Bulletin, EPA/540/MR-93/503.

EPA, 1993. Blogenesis Soll Washing Technology, EPA RREL, series includes Demonstration Bulletin, EPA/540/MR-93/510; innovative Technology Evaluation Report, EPA/540/R-93/510; and

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EPA, 1997. Best Management Practices (BMPs) for Soil Treatment Technologies: Suggested Operational Guidelines to Prevent Cross-media Transfer of Contaminants During Clean-UP Activities, EPA OSWER, EFA/530/R-97/007.

Federal Remediation Technologies Roundtable, 1995. Remediation Case Studies: Thermal Desorption, Soil Washing, and In Situ Vitrification, EPA/542/R-95/005.

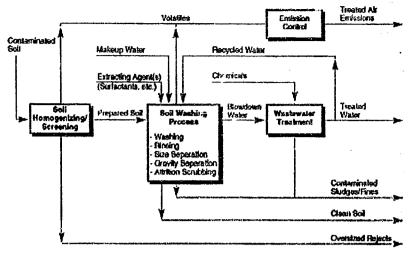
 Soil Washing at the King of Prussia Technical Corporation Superfund Site, Winslow Township, New Jersey

Raghavan, R., D.H. Dietz, and E. Coles, 1988. Cleaning Excuvated Soil Using Extraction Agents: A State-of-the-Art Review, EPA Report EPA 600/2-89/034.

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#### Site Information:

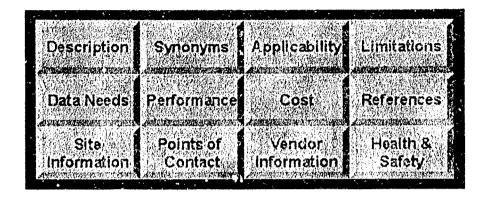
- Ft. Polk, LA
- · Twin Cities AAP, New Brighton, MN
- EPA Demo: Santa Maria, CA
- · Army Saginaw Bay Confined Disposal Facility, MI
- DOE Demo: Clemson Technical Center, SC
- EPA & DOE Demo: Montclair, West Orange & Glen Ridge Sites, NJ
- EPA Demo: Toronto Port Industrial District, Canada
- Army Demo: Sacramento Army Depot, CA
- DOE Demo: Fernald Site, OH
- EPA Demo: Alaska Battery Enterprises Superfund Site, AK
- EPA Demo: MacGil



# 4.24 Solidification/Stabilization

(Ex Situ Soil Remediation Technology)

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Technology	Description	
Soil, Sediment, and Sludge		
3.5 Ex Situ Physical/Chemical Treatment (assuming excavation)		
4.24 Solidification/ Stabilization (Ex Situ)	Contaminants are physically bound or enclosed within a stabilized mass (solidification), or chemical reactions are induced between the stabilizing agent and contaminants to reduce their mobility (stabilization).	

### Description:



Figure 4-24: Typical Ex Situ Solidification/ stabilization Process Flow Diagram

As for in situ solidification/stabilization (S/S) (see <u>Technology Profile No. 4.10</u>), ex situ S/S contaminants are physically bound or enclosed within a stabilized mass (solidification), or chemical reactions are induced between the stabilizing agent and contaminants to reduce their mobility (stabilization). Ex situ S/S, however, typically requires disposal of the resultant materials. Under CERCLA material can be replaced on site.

There are many innovations in the stabilization and solidification technology. Most of the innevations are modifications of proven processes and are directed to encapsulation or immobilizing the harmful constituents and involve processing of the waste or contaminated soil. Nine distinct innovative processes or groups of processes include: (1) bituminization, (2) emulsified asphalt, (3) modified sulfur cement, (4) polyethylene extrusion, (5) pozzolan/Portland cement, (6) radioactive waste solidification, (7) sludge stabilization, (8) soluble phosphates, and (9) vitrification/molten glass.

Typical ex situ S/S is a short- to medium-term technology.

#### > Bituminization

In the bituminization process, wastes are embedded in motion bitumen and encapsulated when the bitumen cools. The process combines heated bitumen and a concentrate of the waste material, usually in slurry form, in a heated extruder containing screws that mix the bitumen and waste. Water is evaporated from the mixture to about 0.5% moisture. The final product is a homogenous mixture of extruded solids and bitumen.

### > Emulsified Asphalt

Asphalt emulsions are very fine droplets of asphalt dispersed in water that are stabilized by chemical emulsifying agents. The emulsions are available as either cationic or anionic emulsions. The emulsified asphalt process involves adding emulsified asphalts having the appropriate charge to hydrophilic liquid or semiliquid wastes at ambient temperature. After mixing, the emulsion breaks, the water in the waste is released, and the organic phase forms a continuous matrix of hydrophobic asphalt around the waste solids. In some cases, additional neutralizing agents, such as lime or gypsum, may be required. After given sufficient time to set and cure, the resulting solid asphalt has the waste uniformly distributed throughout it and is impermeable to water.

### > Modified Sulfur Cement

Modified sulfur cement is a commercially-available thermoplastic material. It is easily melted (127° to 149° C (260° to 300° F)) and then mixed with the waste to form a homogenous molten slurry which is discharged to suitable containers for cooling, storage, and disposal. A variety of common mixing devices, such as, paddle mixers and pug mills, can be used. The relatively low temperatures used limit emissions of sulfur dioxide and hydrogen sulfide to allowable threshold values.

#### > Polyethylene Extrusion

The polyethylene extrusion process involves the mixing of polyethylene binders and dry waste materials using a heated cylinder containing a mixing/transport screw. The heated, homogenous mixture exits the cylinder through an output die into a mold, where it cools and solidifies. Polyethylene's properties produce a very stable, solidified product. The process has been tested on nitrate salt wastes at plant-scale, establishing its viability, and on various other wastes at the bench and pilot scale.

#### > Pozzolan/Portland Cement

Pozzolan/Portland cement process consists primarily of silicates from pozzolanic-based materials like fly ash, kiln dust, pumice, or blast furnace slag and cement-based materials like Portland cement. These materials chemically react with water to form a solid cementious matrix which improves the handling and physical characteristics of the waste. They also raise the pH of the water which may help precipatate and immobilize some heavy metal contaminants. Pozzolanic and cement-based binding agents are typically appropriate for inorganic contaminants. The effectiveness of this binding agent with organic contaminants varies.

### > Radioactive Waste Solidification

In radioactive waste solidification (Grouting/Other) treatment, solidification additives are used to form a uniform and stable matrix to encapsulate radioactive waste materials. Assemblies include pumps for liquids or slurries, conveyors for sludges or solids, storage silos, weigh feeders, piping, mixers and disposal or storage.

#### > Sludge Stabilization

The sludge stabilization process is the addition of a reagent, either slags or cementitious materials, to sludge to transform the material so that the hazardous constituents are in their least mobile or toxic form. Sludges which leach heavy metals or other contaminants are often stabilized to immobilize the hazardous constituents.

#### > Soluble Phosphates

The soluble phosphates process involves the addition of various forms of phosphate and alkali for control of pH as well as for formation of complex metal molecules of low-solubility to immobilize

(insolubilize) the metals over a wide pH range. Unlike most other stabilization processes, soluble phosphate processes do not convert the waste into a hardened, monolithic mass. One application of soluble phosphates and lime is in stabilizing fly ash by immobilizing the lead and cadmium in the ash.

# > Vitrification/Molten Glass

Vitrification, or molten glass, processes are solidification methods that employ heat up to 1,200° C to melt and convert waste materials into glass or other glass and crystalline products. The high temperatures destroy any organic constituents with very few byproducts. Materials, such as heavy metals and radionuclides, are actually incorporated into the glass structure which is, generally, a relatively strong, durable material that is resistant to leaching. In addition to solids, the waste materials can be liquids, wet or dry studges, or combustible materials. Borosilicate and soda-lime are the principal glass formers and provide the basic matrix of the vitrified product.

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Synonyms:

**DSERTS Code:** 

M13 (Vitrification)

N11 (Solidification/Stabilization)

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Applicability:

The target contaminant group for ex situ S/S is inorganics, including radionuclides. Most S/S technologies have limited effectiveness against organics and pesticides, except vitrification which destroys most organic contaminants.

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Limitations:

Factors that may limit the applicability and effectiveness of the process include:

- Environmental conditions may affect the long-term immobilization of contaminants.
- Some processes result in a significant increase in volume (up to double the original volume).
- Certain wastes are incompatible with different processes. Treatability studies are generally required.
- Organics are generally not immobilized.
- Long-term effectiveness has not been demonstrated for many contaminant/process combinations.

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Data Needs:

A detailed discussion of these data elements is provided in Subsection 2.2.1 (Data Requirements for Soil, Sediment, and Sludge). Soil parameters that must be determined include particle size, Atterberg limits, moisture content, metal concentrations, sulfate content, organic content, density, permeability, unconfined compressive strength, leachability, microstructure analysis, and physical and chemical durability.

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Performance Data:

The performance of ex situ S/S is dependent on the type of S/S process used.

DOE has demonstrated the Polyethylene Encapsulation of Radionuclides and Heavy Metals (PERM) process at the bench scale. The process is a waste treatment and stabilization technology for high-level mixed waste. Specific targeted contaminants include radionuclides (e.g., cesium, strontium, and cobalt), and toxic metals (e.g., chromium, lead, and cadmium). Scale-up from bench-scale tests has demonstrated the feasibility to process waste at approximately 2,000 lb/hr. The scale-up feasibility tests have successfully demonstrated the potential to encapsulate at least 60 wt% nitrate salt in polyethylene. Polyethylene waste forms have been demonstrated to exceed Nuclear Regulatory Commission, EPA, and Department of Transportation waste form criteria. Waste forms containing up to several thousand ppm of toxic-metal contaminants have passed the EPA's TCLP.

DOE also demonstrated the arc melter vitrification process, which is capable of melting soil and metals, pyrolizing or oxidizing residual organics, melting structural metals from melted slag (silica and metal oxides), and partitioning transuranic (TRU) waste into slag phase. Durability tests with the resultant slag showed an approximately order of magnitude reduction in leachability when compared with high-level borosilicate glass.

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Cost:

Ex situ solidification/stabilization processes are among the most mature remediation technologies. Representative overall costs from more than a dozen vendors indicate an approximate cost of under \$110 per metric ton (\$100 per ton), including excavation.

# Back to Top

References:

Battelle Memorial Institute, 1995. *ReOpt. V3.1*, by Battelle Memorial Institute for DOE under Contract DE/AC06/76RLO 1830.

Bricka, R.M., et al., 1988. An Evaluation of Stabilization/Solidification of Fluidized Bed Incineration Ash (K048 and K051), USAE-WES Technical Report EL-88-24.

California Base Closure Environmental Committee (CBCEC), 1994. *Treatment Technologies Applications Matrix for Base Closure Activities, Revision 1*, Technology Matching Process Action Team, November, 1994.

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DOE, 1995. *Technology Catalogue, Second Edition*, Office of Environmental Management and Office of Technology Development, DOE/EM-0235.

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EPA, 1989. *Harcon Solidification*, EPA RREL, series includes Technology Evaluation Vol. I, EPA/540/5-89/001a, PB89-158810; Technology Evaluation Vol. II, EPA/540/5-89/001b, PB89-158828; Applications Analysis, EPA/540/A5-89/001; and Technology Demonstration Summary, EPA/540/S5-89/001.

EPA, 1989. Solidtech, Inc. Solidification, EPA RREL, series includes Technology Evaluation Vol. I, EPA/540/5S-89/005a; Technology Evaluation Vol. II, EPA/540/5S-89/005b, PB90-191768; Applications Analysis, EPA/540/A5-89/005; Technology Demonstration Summary, EPA/540/S5-89/005; and Demonstration Bulletin, EPA/540/M5-89/005.

EPA, 1989. Stabilization/Solidification of CERCLA and RCRA Wastes - Physical Tests, Chemical Testing Procedures, Technology Screening and Field Activities, EPA, ORD, Washington, DC, EPA/625/6-89/022.

EPA, 1992. Silicate Technology Corporation Solidification/Stabilization of Organic/Inorganic Contaminants, EPA RREL, Demonstration Bulletin, EPA/540/MR-92/010; Applications Analysis, EPA/540/AR-92/010, PB93-172948.

EPA, 1993. Solidification/Stabilization and Its Application to Waste Materials, Technical Resource Document, EPA, ORD, Washington, DC, EPA/530/R-93/012.

EPA, 1993. Solidification/Stabilization of Organics and Inorganics, Engineering Bulletin, EPA, ORD, Cincinnati, OH, EPA/540/S-92/015.

EPA, 1994. Innovative Site Remediation Technology: Solidification/Stabilization, Vol. 4, EPA OSWER 542/B-94/001.

EPA, 1997. Best Management Practices (BMPs) for Soil Treatment Technologies: Suggested Operational Guidelines to Prevent Cross-media Transfer of Contaminants During Clean-UP Activities, EPA OSWER, EPA/530/R-97/007.

EPA, 1997. Technology Alternatives for the Remediation of Soils Contaminated with As, Cd, Cr, Hg, and Pb, Engineering Bulletin, EPA540/R-97/008.

Federal Remediation Technologies Roundtable, 1997. Remediation Case Studies: Soil Vapor Extraction and Other In Situ Technologies, EPA/542/R-97/009.

• In Situ Enhanced Soil Mixing at the U.S. Department of Energy's Portsmouth Gaseous Diffusion Plant, X-231B Unit, Piketon, Ohio

Federal Remediation Technologies Roundtable, 1998. Remediation Case Studies: Debris and Surface Cleaning Technologies, and Other Miscellaneous Technologies, EPA/542/R-98/017.

Polyethylene Macroencapsulation at Envirocare of Utah, Inc., Salt Lake City, Utah

USAEC, 1997. "Plasma Arc Technology Evaluation" in *Innovative Technology Demonstration*, *Evaluation and Transfer Activities*, FY 96 Annual Report, Report No. SFIM-AEC-ET-CR-97013, pp. 107-110.

Wittle, J.K., et.al., 1995. Graphite Electrode DC Arc Technology Program for Burled Waste Treatment, Electro-Pyrolysis, Inc. Wayne, Penn.

# Back to Top

#### Site Information:

- EPA SITE Demo: Robins AFB, Macon, GA
- EPA SITE Demo: Selma Pressure Treating Selma, CA
- EPA SITE Demo: Portable Equip. Salvage Co. Clackamas, OR
- Navy Demo: Naval Const. Battalion Ctr. Port Hueneme, CA
- Imperial Oil Co./Champion Chemical Co. Superfund Site Morganville, NJ
- Small Arms Range, Naval Air Station Mayport, FL
- Davis-Monthan AFB
- DO! Demo: Salt Lake City Research Center
- DOI Demo: Albany Metallurgy Research Center, OR
- EPA & DOE Demo: Component Development & Integration Facility, MT
- NEL Demo: NAS North Island Installation Restoration (IR) Site 11, CA
- Envirocare of Utah, Inc., Salt Lake City, Utah
- DOE's Portsmouth Gaseous Diffusion Plant, X-231B Unit, Piketon, OH
- Additional site information on the FRTR web site

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#### **Points of Contact:**



**Technology Specific Web Sites:** 



Non-Government Web Sites

Backlo Ton

Vendor Information:

A list of vendors offering Ex Situ Physical/Chemical Soil Treatment is available from the Vendor Information System for Innovative Treatment Technologies (VISITT) developed by U.S. Environmental Protection Agency (EPA).



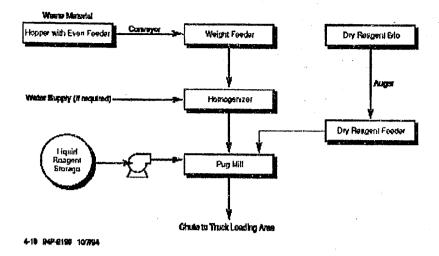
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Health and Safety:



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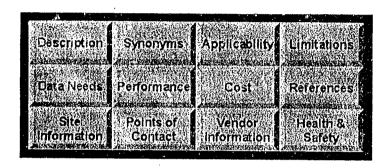
# 4.44 Hydrofracturing

(In Situ Ground Water Remediation Technology)

Previous Section

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Synonym List Next Section



Technology	Description	
Ground Water, Surface Water, and Leachate		
3.10 In Situ Physical/Chemical Treatment		
(enhancement)	Injection of pressurized water through wells cracks low permeability and over-consolidated sediments. Cracks are filled with porous media that serve as substrates for bioremediation or to improve pumping efficiency.	
Decarintions	Hedro Control in 1914 and 1914 and 1914	

# Description:

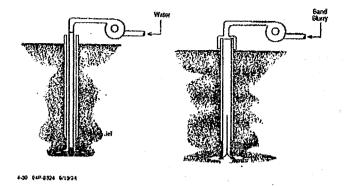


Figure 4-44:
Typical Sequence of
Operations for
Creating Hydraulic
Fractures

Hydrofracturing is a pilot-scale technology in which pressurized water is injected to increase the permeability of consolidated material or relatively impermeable unconsolidated material. Fissures created in the process are filled with a porous medium that can facilitate bioremediation and/or improve extraction efficiency. Fractures promote more uniform delivery of treatment fluids and accelerated extraction of mobilized contaminants. Typical applications are linked with soil vapor extraction, in situ bioremediation, and pump-and-treat systems.

The fracturing process begins with the injection of water into a sealed borehole until the pressure of the water exceeds the overburden pressure and a fracture is created. A slurry composed of a coarse-grained sand and guar gum gel or a similar substitute is then injected as the fracture grows away from the well. After pumping, the sand grains hold the fracture open while an enzyme additive breaks down the viscous fluid. The thinned fluid is pumped from the fracture, forming a permeable subsurface channel suitable for delivery or recovery of a vapor or liquid.

The hydraulic fracturing process can be used in conjunction with soil



vapor extraction technology to enhance recovery. Hydraulically-induced fractures are used to deliver fluids, substrates and nutrients for in situ bioremediation applications.

Back to Tob.

Synonyms:

**DSERTS Code:** F17 (Hydrofracturing - enhancement)

Applicability:

Back to Top :

Hydrofracturing is applicable to a wide range of contaminant groups with no particular target group.

Limitations:

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Factors that may limit the applicability and effectiveness of the process include:

- The technology should not be used in bedrock susceptible to seismic activity.
- Investigation of possible underground utilities, structures, or trapped free product is required.
- The potential exists to open new pathways leading to the unwanted spread of contaminants (e.g., DNAPLs).
- Pockets of low permeability may still remain after using this technology.
- There is an inability to control the final location or size of the fractures that are created.
- Fractures are anticipated to collapse due to over burden pressure.

Dack to Ton

Data Needs:

A detailed discussion of these data elements is provided in Subsection 2.2.2. (Data Requirements for Ground Water, Surface Water, and Leachate).

Dack to Top

Performance Data:

The technology has had widespread use in the petroleum and waterwell construction industries but is an innovative method for remediating hazardous waste sites.

Dack to Ton

Cost:

The cost per fracture is estimated to be \$1,000 to \$1,500, based on creating four to six fractures per day. This cost (including equipment rental, operation, and monitoring) is small compared to the benefits of enhanced remediation and the reduced number of wells needed to complete the remediation. A number of factors affect the estimated costs of creating hydraulic fractures at a site. These factors include physical site conditions such as site accessibility and degree of soil consolidation; degree of soil saturation; and geographical location, which affects availability of services and supplies. The first two factors also affect the effectiveness of hydraulic fracturing.

The costs presented in this analysis are based on conditions found at the Xerox Oak Brook site. A full-scale demonstration was not

conducted for this technology. Because operating costs were not independently monitored during the pilot-scale demonstrations at the Xerox Oak Brook and Dayton sites, all costs presented in this section were provided by Xerox and University of Cincinnati Center Hill.

#### References:

Back to Top

EPA, 1991. Feasibility of Hydraulic Fracturing of Soil to Improve Remedial Actions, EPA/600/S2-91/012.

EPA, 1993. Hydraulic Fracturing Technology, EPA/600/R-93/505.

EPA, 1993. *Hydraulic Fracturing of Contaminated Soil*, series includes Demonstration Bulletin, EPA/540/MR-93/505; Technology Evaluation and Applications Analysis Combined, EPA/540/R-93/505; and Technology Demonstration Summary, EPA/540/SR-93/505.

EPA, 1994. In Situ Remediation Technology Status Report: Hydrofracturing/Pneumatic Fracturing, EPA/542/K-94/005.

EPA, 1997. Analysis of Selected Enhancements for Soil Vapor Extraction, EPA OSWER, EPA/542/R-97/007.

Federal Remediation Technologies Roundtable, 1997. Remediation Case Studies: Soil Vapor Extraction and Other In Situ Technologies, EPA/542/R-97/009.

 Hydraulic and Pneumatic Fracturing at the U.S. Department of Energy's Portsmouth Gaseous Diffusion Plant, Ohio; Department of Defense; and Commercial Sites.

Hubbert, M.K and D.G. Willis, 1957. "Mechanics of Hydraulic Fracturing," *Petroleum Transactions AIME*, Vol. 210, pp. 153 through 168.

Murdoch, L.C., 1990. "A Field Test of Hydraulic Fracturing in Glacial Till," in *Proceedings of the Research Symposium*, Ohio, EPA Report, EPA/600/9-90/006.

Murdoch, L.C., 1993. "Hydraulic Fracturing of Soil During Laboratory Experiments, Part I: Methods and Observations; Part II: Propagation; Part III: Theoretical Analysis", Geotechnique, Vol. 43, No. 2, Institution of Civil Engineers, London, pp. 255 to 287.

University of Cincinnati (UC), 1991. "Work Plan for Hydraulic Fracturing at the Xerox Oak Brook Site in Oak Brook, Illinois".

Wolf, A. and L.C. Murdoch, 1992. "The Effect of Sand-Filled Hydraulic Fractures on Subsurface Air Flow: Summary of SVE Field Tests Conducted at the Center Hill Research Facility", UC Center Hill Facility, Unpublished Report.

## Hack to Top

## Site Information:

- Xerox Facility Oak Brook, IL
- UST site, Dayton, OH
- DOE's Portsmouth Gaseous Diffusion Plant, OH

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# **Points of Contact:**



Technology Specific Web Sites:



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#### **Vendor Information:**

A list of vendors offering In Situ Physical/Chemical Water Treatment is available from the Vendor Information System for Innovative Treatment Technologies (VISITT) developed by U.S. Environmental Protection Agency (EPA).



## Hack to June

## Health and Safety:

To be added

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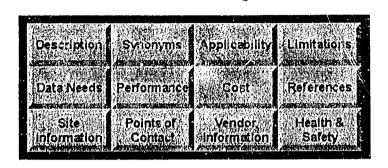
# 4.46 Passive/Reactive Treatment Walls

(In Situ Ground Water Remediation Technology)

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Screen Matrix Table of Contents

Synonym List Next Section



Technology	Description	
Ground Water, Surface Water, and Leachate		
3.10 In Situ Physical/Chemical Treatment		
4.46 Passive/Reactive These barriers allow the passage of water while causing the degradation or removal of contaminants.		

## Description:



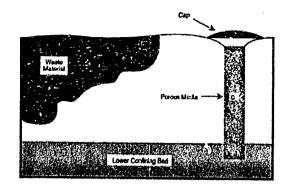
Typical Passive Treatment Wall (Cross-Section)

A permeable reaction wall is installed across the flow path of a contaminant plume, allowing the water portion of the plume to passively move through the wall. These barriers allow the passage of water while prohibiting the movement of contaminants by employing such agents as zero-valent metals, chelators (ligands selected for their specificity for a given metal), sorbents, microbes, and others.

The contaminants will either be degraded or retained in a concentrated form by the barrier material. The wall could provide permanent containment for relatively benign residues or provide a decreased volume of the more toxic contaminants for subsequent treatment.

#### > Funnel and Gate

Modifications to the basic passive treatment walls may involve a funnel-and-gate system or an iron treatment wall. The funnel-and-gate system for in situ treatment of contaminated plumes consists of low hydraulic conductivity (e.g., 1E-6 cm/s) cutoff walls (the funnel) with a gate that contains in situ reaction zones. Ground water primarily flows through high conductivity gaps (the gates). The type of cutoff walls most likely to be used in the current practice are slurry walls or sheet piles. Innovative methods such as deep soil mixing and jet grouting are also being considered for funnel walls.



440 84P-3317 926/34

#### > Iron Treatment Wall

An iron treatment wall consists of iron granules or other iron bearing minerals for the treatment of chlorinated contaminants such as TCE, DCE, and VC. As the iron is oxidized, a chlorine atom is removed from the compound by one or more reductive dechlorination mechanisms, using electrons supplied by the oxidation of iron. The iron granules are dissolved by the process, but the metal disappears so slowly that the remediation barriers can be expected to remain effective for many years, possibly even decades.

Barrier and post-closure monitoring tests are being conducted by the USAF, U.S. Navy, and DOE in field-scale demonstration plots and are being designed for actual contaminated sites. The range of materials available for augmenting existing barrier practice is broad. Two types of barriers have been the focus of initial efforts of this program, i.e., permeable reactive barriers and in-place bioreactors.

Passive treatment walls are generally intended for long-term operation to control migration of contaminants in ground water.

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Synonyms:

Permeable reactive barrier; In place bioreaction; In-situ chemical

DSERTS Code: F16 (Passive Treatment Walls)

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Applicability:

Target contaminant groups for passive treatment walls are VOCs, SVOCs, and inorganics. The technology can be used, but may be less effective, in treating some fuel hydrocarbons.

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Limitations:

Factors that may limit the applicability and effectiveness of the process include:

- Passive treatment walls may lose their reactive capacity, requiring replacement of the reactive medium.
- Passive treatment wall permeability may decrease due to precipitation of metal salts
- Depth and width of barrier.
- Limited to a subsurface lithology that has a continous aquitard at a depth that is within the vertical limits of trenching equipment.
- Volume cost of treatment medium.
- Biological activity or chemical precipitation may limit the permeability of the passive treatment wall.

#### Back to Top

Data Needs:

A detailed discussion of these data elements is provided in <u>Subsection 2.2.2</u>. (Data Requirements for Ground Water, Surface Water, and <u>Leachate</u>).

Data needs include hydraulic gradient; contaminant characteristics (depth, areal extent, type, and concentration); depth to ground water, including range of anticipated fluctuations; depth to impermeable barrier key-in; site stratigraphy; ground water hydrology; water quality, flow rate, and direction; soil permeability; and buffering capacity.

Dackto Top

Performance Data:

Data has been developed by the USAF, the University of Waterloo, and Environmetal Technologies. Inc.

Several full-scale and demonstration scale walls have been installed for remediation of ground water contaminated with chlorinated aliphatic hydrocarbons. These sites include Lowry AFB and Moffett Field NAS. Several more sites are currently being evaluated or have passive treatment walls scheduled for installation.

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Cost:

Complete cost data are still not available because most sites have been demonstration scale and may have been overdesigned for a safety margin. However, costs are decreasing as the price of reactive iron media declines and cost per unit of contaminant removed is a function of the concentrations in ground water.

Dack to Ton.

References:

California Base Closure Environmental Committee (CBCEC), 1994. Treatment Technologies Applications Matrix for Base Closure Activities, Revision 1, Technology Matching Process Action Team, November, 1994.

EPA, 1995. In Situ Remediation Technology Status Report: Treatment Walls, Office of Solid Waste and Emergency Response, EPA/540/K-94/604.

EPA, 1997. Permeable Reactive Subsurface Barriers for the Interception and remediation of Chlorinated Hydrocarbon and Chromium(VI) Plumes in Ground Water, EPA/600/F-97/008.

DOE, 1993. Technical Name: Barriers and Post-Closure Monitoring, Technology Information Profile (Rev. 2), DOE Protech Database, TTP No. AL-1211-25.

DOE, 1994. Technology Catalogue, First Edition. February.

Federal Remediation Technologies Roundtable, 1998. Remediation Case Studies: Innovative Groundwater Treatment Technologies, EPA/542/R-98/015.

- Pump and Treat and Permeable Reactive Barrier to Treat Contaminated Groundwater at the Former Intersil, Inc. Site, Sunnyvale, California
- Permeable Reactive Barrier to Treat Contaminated Groundwater at Moffett Federal Airfield, Mountain View, California
- In Situ Permeable Reactive Barrier for Treatment of Contaminated Groundwater at the U.S. Coast Guard Support Center, Elizabeth City, North Carolina

Hansen, W., et al., 1992. "Barriers and Post-Closure Monitoring", Briefing Chart, Los Alamos National Laboratory, Los Alamos, NM, TTP No. AL-1212-25.

USAF, 1997. Design Guidance for Application of Permeable Barriers to Remediate Dissolved Chlorinated Solvents, prepared by Battelle under contract to Environics Directorate, Armstrong Laboratory.

Vidic, R.D. and F.G. Pohland. "Technology Evaluation Report: Treatment Walls", GWRTAC Series TE-96-01.

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#### Site Information:

- Tonolli Corporation, PA
- Lowry Air Force Base, CO
- Somersworth Sanitary Landfill Site, NH
- Moffett Field NAS, CA
- Brown's Battery Breaking Site, OU 2, PA
- Canadian Forces Base, Canada
- Borden Aquifer, Canada
- · Air Force Demo Hill AFB, UT
- Los Alamos National Laboratory
- Former Intersil, Inc. Site, Sunnyvale, CA
- · Moffett Federal Airfield, Mountain View, CA
- The U.S. Coast Guard Support Center, Elizabeth City, NC
- Additional site information on the FRTR web site

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## **Points of Contact:**



**Technology Specific Web Sites:** 

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# Non-Government Web Sites

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# Vendor Information:

A list of vendors offering In Situ Physical/Chemical Water Treatment is available from the Vendor Information System for innovative Treatment Technologies (VISITT) developed by U.S. Environmental Protection Agency (EPA).



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Health and Safety:



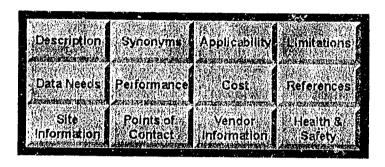
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# 4.57 Deep Well Injection

(Ground Water Containment Remediation Technology)

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Technology	Description	
Ground Water, Surface Water, and Leachate		
3.13 Containment		
Injection	Deep well injection is a liquid waste disposal technology. This alternative uses injection wells to place treated or untreated liquid waste into geologic formations that have no potential to allow migration of contaminants into potential potable water aquifers.	

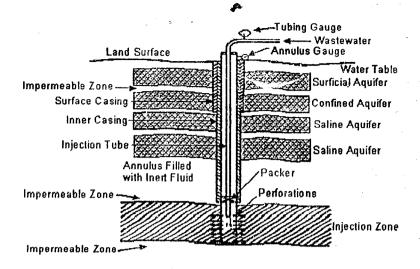
#### Description:



Figure 4-57: Typical Deep Well Injection System

A typical injection well consists of concentric pipes, which extend several thousand feet down from the surface level into highly saline, permeable injection zones that are confined vertically by impermeable strata. The outermost pipe or surface casing, extends below the base of any underground sources of drinking water (USDW) and is cemented back to the surface to prevent contamination of the USDW. Directly inside the surface casing is a long string casing that extends to and sometimes into the injection zone. This casing is filled in with cement all the way back to the surface in order to seal off the injected waste from the formations above the injection zone back to the surface. The casing provides a scal between the wastes in the injection zone and the upper formations. The waste is injected through the injection tubing inside the long string easing either through perforations in the long string or in the open hole below the bottom of the long string. The space between the string casing and the injection tube, called the annulus, is filled with an inert, pressurized fluid, and is sealed at the bottom by a removable packer preventing injected wastewater from backing up into the annulus.

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Synonyms:

Subsurface injection, Underground injection, Class I injection wells.

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Applicability:

The target contaminant groups for deep well injection are VOCs, SVOCs, fuels, explosives, and pesticides. However, existing permitted deep well injection facilities are limited to a narrow range of specific wastes. Success at expanding existing permits to manage hazardous wastes seems unlikely.

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Limitations:

Factors that may limit the applicability and effectiveness of these processes include:

- Injection will not be used for hazardous waste disposal in any areas where seismic activity could potentially occur.
- Injected wastes must be compatible with the mechanical components of the injection well system and the natural formation water. The waste generator may be required to perform physical, chemical, biological, or thermal treatment for removal of various contaminants or constituents from the waste to modify the physical and chemical character of the waste to assure compatibility.
- High concentrations of suspended solids (typically >2 ppm) can lead to plugging of the injection interval.
- Corrosive media may react with the injection well components, with injection zone formation, or with confining strata with very undesirable results. Wastes should be neutralized.
- High iron concentrations may result in fouling when conditions alter the valence state and convert soluble to insoluble species.
- Organic carbon may serve as an energy source for indigenous or injected bacteria resulting in rapid population growth and subsequent fouling.
- Waste streams containing organic contaminants above their solubility limits may require pretreatment before injection into a well.
- Site assessment and aquifer characterization are required to determine suitability of site for wastewater injection.
- Extensive assessments must be completed prior to receiving approval from regulatory authority.

TDack to Top 6

Data Needs:

A detailed discussion of data elements is provided in <u>Subsection 2.2.2</u> (Data Requirements for Ground Water, Surface Water, and Leachate).

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Performance Data:

Injection wells have been used for the disposal of industrial and hazardous wastes since the 1950s, so the equipment and methodology are readily available and well known; however the use of them today is continuing under very strict regulatory control.

Dack to Top-

Cost:

NA

References:

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EPA, 1985. Report to Congress on Injection of Hazardous Waste,

EPA, Office of Drinking Water, EPA/9-85-003.

Reeder et al., 1977. Review and Assessment of Deep Well Injection

of Hazardous Waste, Volume I, EPA/600/2-77/029a.

Warner and J.H.Lehr, 1977. An Introduction to the Technology of

Subsurface Wastewater Injection, EPA/600/2-77/240.

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Site Information:

• DOI Demo Bureau of Mines Tuscaloosa Research Center, AL

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**Points of Contact:** 

General/FRTR Agency Contacts

Technology Specific Web Sites:

NA

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**Vendor Information:** 

A list of vendors offering Water Containment Treatment is available from the Vendor Information System for Innovative Treatment Technologies (VISITT) developed by U.S. Environmental Protection Agency (EPA).



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Health and Safety:

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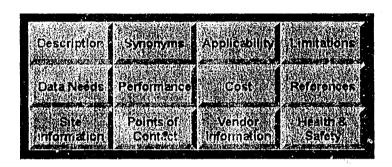
# APPENDIX C

Theoretically Feasible Technologies

# 4.48 Constructed Wetlands

(Ex Situ Ground Water Remediation Technology)

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Section Page Matrix Contents List Section



Technology	Description	
Ground Water, Surface Water, and Leachate		
3.11 Ex Situ Biological Treatment (assuming pumping)		
Wetlands	The constructed wetlands-based treatment technology uses natural geochemical and biological processes inherent in an artificial wetland ecosystem to accumulate and remove metals, explosives, and other contaminants from influent waters. The process can use a filtration or degradation process.	

Description:

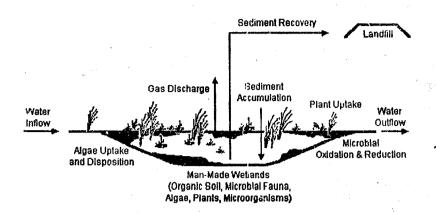


Figure 4-48: Typical Constructed Wetlands System

Although the technology incorporates principal components of wetland ecosystems, including organic soils, microbial fauna, algae, and vascular plants, microbial activity is responsible for most of the remediation.

Influent waters with high metal concentrations and low pH flow through the aerobic and anaerobic zones of the wetland ecosystem. Metals are removed through ion exchange, adsorption, absorption, and precipitation with geochemical and microbial oxidation and reduction. Ion exchange occurs as metals in the water contact humic or other organic substances in the wetland. Wetlands constructed for this purpose often have little or no soil instead they have straw, manure or compost. Oxidation and reduction reactions catalyzed by bacteria that occur in the aerobic and anaerobic zones, respectively, play a major role in precipitating metals as hydroxides and sulfides. Precipitated and adsorbed metals settle in quiescent ponds or are filtered out as water percolates through the medium or the plants.

Influent water with explosive residues or other contaminants flows through and beneath the gravel surface of a gravel-based wetland. The



wetland, using emergent plants, is a coupled anaerobic-aerobic system. The anaerobic cell uses plants in concert with natural microbes to degrade the contaminant. The aerobic, also known as the reciprocating, cell further improves water quality through continued exposure to the plants and the movement of water between cell compartments.

Wetland treatment is a long-term technology intended to operate continously for years.

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Synonyms:

NA

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Applicability:

Constructed wetlands have most commonly been used in wastewater treatment for controlling organic matter; nutrients, such as nitrogen and phosphorus; and suspended sediments. The wetlands process is also suitable for controlling trace metals, and other toxic materials. Additionally, the treatment has been used to treat acid mine drainage generated by metal or coal mining activities. These wastes typically contain high metals concentrations and are acidic. The process can be adapted to treat neutral and basic tailings solutions.

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Limitations:

The wetlands remediation technology must be adjusted to account for differences in geology, terrain, trace metal composition, and climate in the metal mining regions of the western United States.

The following factors may limit the applicability and effectiveness of the process:

- The long-term effectiveness of constructed wetlands is not well known. Wetland aging may be a problem which may contribute to a decrease in contaminant removal rates over time.
- The cost of building an artificial wetland varies considerably from project and may not be financially viable for many sites.

Back to Ton

Data Needs:

Performance Data:

A detailed discussion of data elements is provided in <u>Subsection 2.2.2</u> (Data Requirements for Ground Water, Surface Water, and Leachate). This technology was accepted into the Emerging Technology Program in 1988; the project was completed in 1991. The purpose of the project was to build, operate, monitor, and assess the effectiveness of a constructed wetlands in treating a portion of acid mine drainage from the Big Five Tunnel site near Idaho Springs, Colorado. The Final Report (EPA/540/R-93/523) is available from NTIS (Order No. PB93-233914). The Summary (EPA/540/SR-93/523) and Bulletin

(EPA/540/F-92/001) are available from EPA.

Study results indicated that heavy metal removal efficiency can approach the removal efficiency of chemical precipitation treatment plants. Some of the optimum results from the 3 years of operation are

#### listed below.

• pH was raised from 2.9 to 6.5.

• Dissolved aluminum, cadmium, chromium, copper, and zinc concentrations were reduced by 99 percent or more.

• Iron was reduced by 99 percent.

• Lead was reduced by 94 percent or more.

Nickel was reduced by 84 percent or more.

 Manganese removal was relatively low, with reduction between 9 and 44 percent.

 Biotoxicity to fathead minnows and water fleas was reduced by factors of 4 to 20.

Because wetland removal processes are primarily microbial, the technology can be developed with traditional process engineering approaches. Laboratory studies can indicate whether remediation is possible, while bench-scale experiments can determine the proper loading and reactor design. Using this approach, five laboratory proof-of-principle studies and three bench-scale studies have been performed, and at least four successful demonstration reactors have been built to remove heavy metals from different types of water.

A final project goal was to develop a manual that discusses design and operating criteria for constructing a full-scale wetland to treat acid mine discharges. The "Wetland Designs for Mining Operations" manual is available from the National Technical Information Service.

Based on the results from the SITE Emerging Technology Program, this technology was selected to participate in the SITE Demonstration Technology Program. Under the Demonstration Program, EPA is evaluating the effectiveness of biogeochemical processes at the Burleigh Tunnel mine site, near Silver Plume, Colorado. Treatment of Burleigh Tunnel discharge is part of the remedy for the Clear Creek/Central City Superfund site. Construction of a pilot-scale system began in summer 1993 and was completed in October 1993. For more information on this project, refer to the Colorado Department of Health profile in the Demonstration Program section (ongoing projects).

The USAEC is demonstrating a gravel based wetland system at Milan Army Ammunition Plant through the ESTCP. The gravel-based system has been effective at degrading TNT and RDX, with a total nitrobody concentration of 10,000 ppb. Analyses indicate degradation due to the rise and fall of daughter products. TNT is reduced to less than 2 ppb. The demonstration had been operational since June 1996.

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Cost:

Studies at Milan Army Ammunition Plant conducted by USAEC and Tennessee Valley Authority (TVA) indicate that plants with nitroreductase, in concert with microbes, can degrade explosive residues. It is estimated that amortizing the capital costs of wetland treatment over a 10 year period results in a cost of \$1.36/Kgal; over a

30 years period, the cost is \$0.45/Kgal.

Back to Top :

References:

EPA, 1993. Constructed Wetlands-Based Treatment, EPA/540/R-

93/523.

EPA, 1995. Colorado School of Mines profile in the Emerging

Technology Program.

Lefave, J.P., 1997. Constructed Wetlands for Treatment of NPS

Pollution, NFESC.

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#### Site Information:

Burleigh Tunnel Silver Plume, CO

• EPA Demo Burleigh Tunnel, CO

Big Five Tunnel, CO

Bear Creek, Oak Ridge, TN

• Naval Amphibious Base in Little Creek, Virginia

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**Points of Contact:** 



**Technology Specific Web Sites:** 



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#### Vendor Information:

A list of vendors offering Ex Situ Biological Water Treatment is available from the Vendor Information System for Innovative Treatment Technologies (VISITT) developed by U.S. Environmental Protection Agency (EPA).



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# Health and Safety:

To be added





# 4.49 Adsorption/Absorption

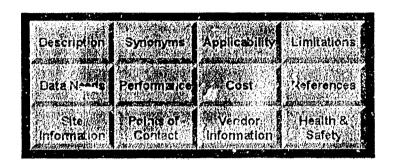
(Ex Situ Ground Water Remediation Technology)

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Synonym List

Next :



Technology	Description	
Ground Water, Surface Water, and Leachate		
3.12 Ex Situ Physical/Chemical Treatment		
	In liquid adsorption, solutes concentrate at the surface of a sorbent, thereby reducing their concentration in the bulk liquid phase.	

# Description:

fro in

Figure 4-49 Typical Adsorption/ Absorption System Adsorption mechanisms are generally categorized as either physical adsorption, chemisorption, or electrostatic adsorption. Neak molecular forces, such as Van der Waals forces, provide the driving force for physical adsorption, while a chemical reaction forms a chemical bond between the compound and the surface of the solid in chemisorption. Electrostatic adsorption involves the adsorption of ions through Coulombic forces, and is normally referred to as ion exchange, which is addressed separately in the ion exchange modules. In liquids, interactions between the solute and the solvent also play an important role in establishing the degree of adsorption.

The most common adsorbent is granulated activated carbon (GAC) (Technology Profile No. 4.51). Other natural and synthetic adsorbents include: activated alumina, forage sponge, lignin adsorption, sorption clays, and synthetic resins.

#### > Activated Alumina

Activated alumina is a filter media made by treating aluminum ore so that it becomes porous and highly adsorptive. Activated alumina will remove variety of contaminants, including excessive fluoride, arsenic, and selenium. The medium requires periodic cleaning with an appropriate regenerant such as alum or acid in order to remain

effective.

# > Forage Sponge

Forage sponge is an open-celled cellulose sponge incorporating an amine-containing chelating polymer that selectively absorbs dissolved heavy metals. The polymer is intimately bonded to the cellulose so as to minimize physical separation from the supporting matrix. The functional groups in the polymer (i.e. amine and carboxyl groups) provide selective affinity for heavy metals in both cationic and anionic states, preferentially forming complexes with transition-group heavy metals.

# ➤ Lignin Adsorption/Sorptive Clay

Lignin adsorption/sorptive clays are used to treat aqueous waste streams with organic, inorganic and heavy metals contamination. The waste stream is treated due to the molecular adhesion of the contaminants to an adsorptive surface.

# > Synthetic Resins

Synthetic resins are more expensive than GAC, but can be designed to achieve higher degrees of selectivity and adsorption capacity for certain compounds than activated carbon. Resins are typically regenerated using acids, bases, or organic sc vents, instead of thermal methods, so they are better suited for thermally unstable compounds such as explosives, and are resistant to deactivation due to the adsorption of dissolved solids. Additionally, resins tend to be more resistant to abrasion than activated carbon, increasing their service life.

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Synonyms:

Liquid phase adsorption.

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Applicability:

The target contaminants groups for adsorption/absorption processes are most organic contaminants and selected inorganic contaminants from liquid and gas streams. Activated alumina can remove fluoride and heavy metals. The forager sponge is specifically used to remove heavy metals. Lignin adsorption/sorptive clays treat organic, inorganic and heavy metals contamination within aqueous waste streams. Synthetic resins are better suited for thermally unstable compounds such as explosives than GAC, due to the resins' non-thermal regeneration requirements.

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Limitations:

Factors that may limit the applicability and effectiveness of these processes include:

- Water-soluble compounds and small molecules are not adsorbed well.
- Costs are high if used as the primary treatment on waste streams with high contaminant concentration lev is.
- Not applicable to sites having high levels of oily substances.
- Not practical where the content of the absorbable hazardous substance is so high that very frequent replacement of the absorbent unit is necessary.
- Contaminated media often require treatment/disposal as hazardous wastes, if they can't be regenerated.

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Data Needs:

A detailed discussion of data elements is provided in Subsection 2.2.2 (Data Requirements for Ground Water, Surface Water, and Leachate).

# Back to Tup

Performance Data:

Adsorption/absorption processes have a long history of use as treatment for municipal, industrial, and hazardous waste streams. The concepts, theory, and engineering aspects of the technologies are well developed. They are proven technologies with documented performance data.

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Cost:

The cost to treat heavy metal contaminated ground water over a one year period with the Forager Sponge technology is estimated at \$340/1000 gallons, assuming the sponges are not regenerated and are replaced upon saturation; or \$238/1000 gallons, assuming the sponges are regenerated twice providing for three useful treatment cycles. Costs for other adsorbent processes are not available.

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References:

Battelle Memorial Institute, 1995. "ReOpt V3.1 User Documentation", for DOE under contract DE/AC06/76RLO 1830.

Rainer, N., 1995. "Forager Sponge, Tech. logy Description", Dynaphone, Inc.

Water Quality Association, 1994. "Treating the Water We Drink, When and Where We Drink It.", WaterReview Technical Brief, Vol. 9, No. 4.

#### Back to Top

#### Site Information:

- EPA Demo National Lead Industry, NJ
- Pease Air Force Base, Newington, NH
  Superfund Site: Western Processing, WA

- Superfund Site: Rocky Mountain Arsenal, OU 17, CO
- Superfund Site: Vally Wood Preserving, Inc., Turlock, CA
- Superfund Site: Motor Wheel, Lansing TWP, MI
- Superfund Site: Perham Arsenic, Perham, MN
- Additional site information on the FRTR web site

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### **Points of Contact:**



# **Technology Specific Web Site:**



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## Vendor Information:

A list of vendors offering Ex Situ Physical/Chemical Water Treatment is available from the Vendor Information System for Innovative Treatment Technologies (VISITT) developed by U.S. Environmental Protection Agency (EPA).



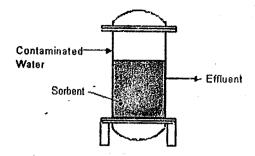
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# Health and Safety:



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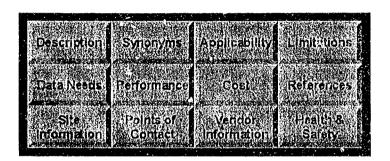
# 4.52 Ion Exchange

(Ex Situ Ground Water Remediation Technology)

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Synonym List Next Section



Technology	Description	
Ground Water, Surface Water, and Leachate		
3.12 Ex Situ Physical/Chemical Treatment (assuming pumping)		
	Ion exchange removes ions from the aqueous phase by exchange with counter ions on the exchange medium.	

Description:



Typical Ion Exchange and Adsorption Equipment Diagram

Ion exchange removes ions from the aqueous phase by the exchange of cations or anions between the contaminants and the exchange medium. Ion exchange materials may consist of resins made from synthetic organic materials that contain ionic functional groups to which exchangeable ions are attached. They also may be inorganic and natural polymeric materials. After the resin capacity has been exhausted, resins can be regenerated for re-use.

The duration of ion exchange technology is typically short- to medium-term depending on the factors discussed in Data Needs.

Synonyms:

NA

Élack (d.To))

Dack to Top'

Applicability:

Ion exchange can remove dissolved metals and radionuclides from aqueous solutions. Other compounds that have been treated include nitrate, ammonia nitrogen, and silicate.

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Principle of infeed-bed ion exchanger (a) Service period. (b) Backwash period. (c) Simultaneous regeneration. (illinois Water Treatment Co.)

Bosros: Chemical Engineer's Handbook, Perry & Clatton (5th Edition

4-46 E4P-5118 9/1992

#### Limitations:

Factors that may affect the applicability and effectiveness of this process include:

- Oil and grease in the ground water may clog the exchange resin.
- Suspended solids content greater than 10 ppm may cause resin blinding.
- The pH of the influent water may affect the ion exchange resin selection.
- Oxidants in ground water may damage the ion exchange resin.
- Wastewater is generated during the regeneration step and will require additional treatment and disposal.

#### Back to Top.

### Data Needs:

A detailed discussion of these data elements is provided in <u>Subsection</u> 2.2.2 (Data Requirements for Ground Water, Surface Water, and <u>Leachate</u>).

Factors affecting the design of an ion exchange system include the presence of oil and grease, contaminant concentration, exchange capacity of the resin, suspended solids, metals, oxidant content, inorganic ions in ground water; and pH of the ground water.

#### Back to Top?

#### Performance Data:

DOE has developed compact processing units (CPUs), or "modular waste treatment units," which are relatively small mobile equipment modules. They perform unit chemical process operations. The CPUs allow rapid deployment of technologies for the treatment of radioactive wastes in underground storage tanks. The modules would be manufactured off-site by commercial vendors and moved into place using trucks or special transports. The concept of having standardized modules is based on the notion that various radioactive waste treatment subsystems could be standardized to match the CPU hardware package, leading to more rapid, cost-effective deployment. The cost benefits are realized even when multiple units are deployed to achieve greater processing rates. The modular design concept will also allow for reuse of CPU components for different unit processes or process deployments.

The ion-exchange CPU will pump undiluted liquid tank waste from an underground storage tank or receive liquid waste from a waste retrieval system for treatment. DOE Northwest Laboratories developed the CPU concept in FY91. Development of a cesium ion-exchange CPU technology is scheduled for 1996. A radioactive waste treatment demonstration is scheduled for FY97.

Another DOE technology, the resorcinol-formaldehyde ion exchange (ReFIX) resin, is being developed for prototype demonstration at the Hanford site in FY97. ReFIX resin is applicable to high-level wastestreams containing cesium-supernate salt solutions.

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Cost:

The cost for a typical ion exchange system ranges from \$0.08 to \$0.21 per 1,000 liters (\$0.30 to \$0.80 per 1,000 gallons) treated. Key cost factors include:

- Pretreatment requirements.
- Discharge requirements and resin utilization.
- Regenerant used and efficiency.

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References:

California Base Closure Environmental Committee (CBCEC), 1994. Treatment Technologies Applications Matrix for Base Closure Activities, Revision 1, Technology Matching Process Action Team, November, 1994.

DOE, 1993. Technology Name: Cesium Removal by Compact Processing Units for Radioactive Waste Treatment, Technology Information Profile (Rev. 2) for ProTech, DOE ProTech Database, TTP Reference No.: RL-321221.

DOE, 1993. Technology Name: Resorcinol-Formaldehyde Ion Exchange Resin for Elutable Ion Exchange in the Compact Portable Units (CPUs) Proposed at Hanford, Technology Information Profile (Rev. 2) for ProTech, DOE ProTech Database, TTP Reference No.: SR-1320-02.

DOE, 1994. Technology Catalogue, First Edition. February.

EPA, 1990. Innovative and Alternative Technology Assessment Manual, EPA, Office of Water Program Operations, EPA/430/9-78/009.

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### **Site Information:**

- EPA Demo Kerr-McGee Chemical Corp., WI
- Army Demo USACE-WES, MS
- Additional site information on the FRTR web site

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**Points of Contact:** 



**Technology Specific Web Sites:** 





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# Vendor Information:

A list of vendors offering Ex Situ Physical/Chemical Water Treatment is available from the Vendor Information System for Innovative Treatment Technologies (VISITT) developed by U.S. Environmental Protection Agency (EPA).



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# Health and Safety:



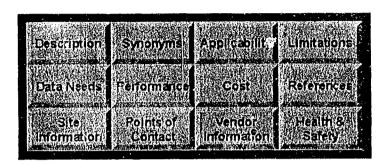
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# 4.53 Precipitation/Coagulation/Flocculation

(Ex Situ Ground Water Remediation Technology)

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Section Page Matrix Contents List Section



Technology	Description	
Ground Water, Surface Water, and Leachate		
3.12 Ex Situ Physical/Chemical Treatment (assuming pumping)		
4.53 Precipitation/ Coagulation/ Flocculation	This process transforms dissolved contaminants into an insoluble solid, facilitating the contaminant's subsequent removal from the liquid phase by sedimentation or filtration. The process usually uses pH adjustment, addition of a chemical precipitant, and flocculation.	

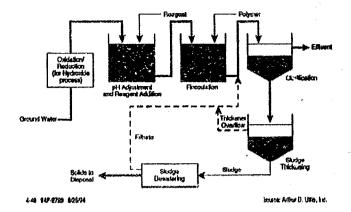
#### Description:



Typical Metals
Precipitation Process

Precipitation of metals has long been the primary method of treating metal-laden industrial wastewaters. As a result of the success of metals precipitation in such applications, the technology is being considered and selected for use in remediating ground water containing heavy metals, including their radioactive isotopes. In ground water treatment applications, the metal precipitation process is often used as a pretreatment for other treatment technologies (such as chemical oxidation or air stripping) where the presence of metals would interfere with the other treatment processes.

Metals precipitation from contaminated water involves the conversion of soluble heavy metal salts to insoluble salts that will precipitate. The precipitate can then be removed from the treated water by physical methods such as clarification (settling) and/or filtration. The process usually uses pH adjustment, addition of a chemical precipitant, and flocculation. Typically, metals precipitate from the solution as hydroxides, sulfides, or carbonates. The solubilities of the specific metal contaminants and the required cleanup standards will dictate the process used. In some cases, process design will allow for the generation of sludges that can be sent to recyclers for metal recovery.



# > Coagulants and Flocculation

In the precipitation process, chemical precipitants, coagulants, and flocculantation are used to increase particle size through aggregation. The precipitation process can generate very fine particles that are held in suspension by electrostatic surface charges. These charges cause clouds of counter-ions to form around the particles, giving rise to repulsive forces that prevent aggregation and reduce the effectiveness of subsequent solid-liquid separation processes. Therefore, chemical coagulants are often added to overcome the repulsive forces of the particles. The three main types of coagulants are inorganic electrolytes (such as alum, lime, ferric chloride, and ferrous sulfate), organic polymers, and synthetic polyelectrolytes with anionic or cationic functional groups. The addition of coagulants is followed by low-sheer mixing in a flocculator to promote contact between the particles, allowing particle growth through the sedimentation phenomenon called flocculant settling.

Flocculant settling refers to a rather dilute suspension of particles that coalesce, or flocculate, during the sedimentation operation. As coalescence or flocculation occurs, the particles increase in mass and settle at a faster rate. The amount of flocculation that occurs depends on the opportunity for contact, which varies with the overflow rate, the depth of the basin, the velocity gradients in the system, the concentration of particles, and the range of particles sizes. The effects of these variables can only be accomplished by sedimentation tests.

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Synonyms:

NA.

Applicability:

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Precipitation is used mainly to convert dissolved ionic species into solid-phase particulates that can be removed from the aqueous phase by coagulation and filtration. Remedial application of this technology usually involve removal of dissolved toxic metals and radionuclides. Depending on the process design, sludges may be amenable to metal recovery.

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Limitations:

Disadvantages of metals precipitation may include:

- As with any pump and treat process, if the source of contamination is not removed (as in metals absorbed to soil), treatment of the ground water may be superfluous.
- The presence of multiple metal species may lead to removal difficulties as a result of amphoteric natures of different compounds (i.e., optimization on one metal species may prevent removal of another).
- As discharge standards become more stringent, further treatment may be required.
- Metal hydroxide sludges must pass TCLP prior to land

disposal.

 Soluble hexavalent chrome requires extra treatment prior to coagulaation and flocculation.

 Reagent addition must be carefully controlled to preclude unacceptable concentrations in treatment effluent.

• Efficacy of the system relies on adequate solids separation techniques (e.g., clarification, flocculation, and/or filtration).

Process may generate toxic sludge requiring proper disposal.

 Process can be costly, depending on reagents used, required system controls, and required operator involvement in system operation.

 Dissolved salts are added to the treated water as a result of pH adjustment.

 Polymer may need to be added to the water to achieve adequate settling of solids.

• Treated water will often require pH adjustment.

 Metals held in solution by complexing agents (e.g., cyanide or EDTA) are difficult to precipitate.

#### Back to Top:

#### Data Needs:

A detailed discussion of these data elements is provided in <u>Subsection</u> 2.2.2 (Data Requirements for Ground Water, Surface Water, and Leachate).

Bench-scale treatability tests should be conducted to determine operating parameters and characteristics [i.e., reagent type and dosage, optimum pH, retention time, flow rate, temperature, mixing requirements, flocculent (polymer) selection, suspended solids, precipitate settling and filtration rates, and sludge volume and characteristics].

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#### Performance Data:

Precipitation of heavy metals as the metal hydroxides or sulfides has been practiced as the prime method of treatment for heavy metals in industrial wastewater for many years. More recently, precipitation (usually as the metal hydroxides) has been used in the electronics and electroplating industries as a pretreatment technology for wastewater discharge to a publicly owned treatment works (POTW). Metals precipitation is widely used to meet NPDES requirements for the treatment of heavy metal-containing wastewaters.

Because of its success in meeting requirements for discharge of treated wastewater, metals precipitation is recognized as a proven process for use in remedial activities such as ground water treatment. Precipitation (combined with sedimentation, and/or flocculation and filtration) is becoming the most widely selected means for heavy metals removal from ground water in pump and treat operations.

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Cost:

The primary capital cost factor is design flow rate. Capital costs for 75- and 250-liters-per-minute (20-gpm and 65-gpm) packaged metals precipitation systems are approximately \$85,000 and \$115,000, respectively.

The primary factors affecting operating costs are labor and chemical costs. Operating costs (excluding sludge disposal) are typically in a range from \$0.08 to \$0.18 per 1,000 liters (\$0.30 to \$0.70 per 1,000 gallons) of ground water containing up to 100 mg/L of metals.

For budgetary purposes, sludge disposal may be estimated to increase operating costs by approximately \$0.13 per 1,000 liters (\$0.50 per 1,000 gallons) of ground water treated. Actual sludge disposal costs (including fixation and transportation) have been estimated at approximately \$330 per metric ton (\$300 per ton) of sludge.

Costs for performing a laboratory treatability study for metals precipitation may range from \$5,000 to \$20,000. Depending on the degree of uncertainty or other requirements, a pilot or field demonstration may be needed. Associated costs may range from \$50,000 to \$250,000 (depending on scale, analytical requirements, and duration).

Additional cost information can be found in the Hazardous, Toxic, and Radioactive Wastes (HTRW) Historical Cost Analysis System (HCAS) developed by Environmental Historical Cost Committee of Interagency Cost Estimation Group.

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References:

Balaso, C.A., et al., 1986. "Soluble Sulfide Precipitation Study", Arthur D. Little, Inc., Final Report to USATHAMA, Report No. AMXTH-TE-CR-87106.

Battelle Memorial Institute, 1995. "ReOpt. V3.1", by Battelle Memorial Institute for DOE under Contract DE/AC06/76RLO 1830.

Bricka, R. Mark, 1988. "Investigation and Evaluation of the Performance of Solidified Cellulose and Starch Xanthate Heavy Metal Sludges", USACE-WES Technical Report EL-88-5.

EPA, 1980. Control and Treatment Technology for the Metal Finishing Industry: Sulfide Precipitation, EPA/625/8-80/003.

EPA, 1990. Innovative and Alternative Technology Assessment Manual, EPA, Office of Water Program Operations, EPA/430/9-78/009.

Federal Remediation Technologies Roundtable, 1998. Remediation Case Studies: Innovative Groundwater Treatment Technologies, EPA/542/R-98/015.

 Coagulation/Flocculation/Dissolved Air Flotation and Oleofiltration<sup>TM</sup> at the Coastal Systems Station, AOC 1, Panama City, Florida

NEESA, 1993. Precipitation of Metals from Ground Water. NEESA Document Number 20.2-051.6, Naval Energy and Environmental Support Activity, Port Hueneme, CA.

Tchobanoglous, G. and F.L. Burton, 1991. "Wastewater Engineering - Treatment, Disposal and Reuse," Third Edition. Metcalf & Eddy, Inc.

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# Site Information:

- EPA Demo Palangana Uranium Mine Site, TX
- Coakley Landfill New Hampshire
- Stringfellow Acid Pit Site CA
- Winthrop Landfill Winthrop, ME
- EPA Removal Action Crown Plating, MO
- Coastal Systems Station, AOC 1, Panama City, FL
- Additional site information on the FRTR web site

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# **Points of Contact:**



Technology Specific Web Sites:



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#### Vendor Information:

A list of vendors offering En Situ Physical/Chemical Water Treatment is available from the Vendor Information System for Innovative Treatment Technologies (VISITT) developed by U.S. Environmental Protection Agency (EPA).



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# 4.40 Directional Wells

(In Situ Ground Water Remediation Technology)

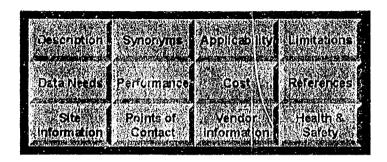
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Technology

Description

Ground Water, Surface Water, and Leachate

3.10 In Situ Physical/Chemical Treatment

4.40 Directional Wells Drilling techniques are used to position wells horizontally, or at an angle, to reach contaminants not accessible by direct vertical drilling.

Description:



Figure 4-40:

Horizontal Wells

Drilling techniques are used to position wells horizontally, or at an angle, to reach contaminants not accessible by direct vertical drilling. Directional drilling may be used to enhance other in-situ or in-well technologies such as ground water pumping, bioventing, SVE, soil flushing, and in-well air stripping.

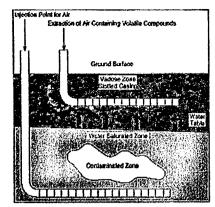
Typical Diagram of In Hardware used for directional boring includes wireline coring rigs, Situ Air Stripping with hydraulic thrust systems, electric cone penetrometers, steering tracking hardware, sonic drilling, and push coring systems. Hydraulically activated thrust equipment capable of exerting more than 40 tons of thrust is used to push the directional boring heads into the earth. Directional control is obtained by proper positioning of the face of the nonsymmetric boring head. Slow rotation of the boring head will cut and compact the geologic material into the borehole wall. Thrusting a boring head that is not rotating wil have a directional change. The machinery is capable of initiating a borehole. steering down to a desired horizontal depth, continuing at that depth, and then steering back to the surface at a downrange location.

Synonyms:

Horizontal Wells.

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146 147-1431 8269

Applicability:

Directional well technology is applicable to the complete range of contaminant groups with no particular target group. It is particularly useful when existing structures interfere with placement of vertical wells.

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Limitations:

Factors that may limit the applicability and effectiveness of the process include:

- The potential exists for the wells to collapse.
- Specialized equipment is required.
- Wells are difficult to position precisely.
- Installation of horizontal wells is typically costly.
- Currently, the technology is limited to depths of less than 50 feet.

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**Data Needs:** 

A detailed discussion of these data elements is provided in <u>Subsection 2.2.2</u>. (Data Requirements for Ground Water, Surface Water, and Leachate).

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Performance Data:

Testing was performed as part of the Mixed Waste Landfill Integrated Demonstration at Sandia National Laboratories, Albuquerque, NM. Several directional holes were drilled; a depth of 12 meters (40 feet) was achieved with a maximum horizontal extent of 174 meters (570 feet).

A DOE field demonstration at the Savannah River site was performed in FY90 for in situ air stripping (ISAS), a mass transfer process that uses horizontal injection and vacuum extraction wells to remediate sites contaminated with VOCs within the vadose zone and soil/ground water in the saturated zone. Air is injected into the saturated zone through horizontal injection wells placed below the water table. As the air passes through the contaminant plume, it volatilizes the chemical constituents. This process performs best in homogeneous soil conditions, while heterogeneities such as formations, fractures. clay layers, and partial clay lenses hinder performance. Clay layers often have high contaminant concentrations, while stratigraphy can cause preferential flow paths and limit the process efficiency. ISAS has been shown to be effective when some interbedded, thin, and/or discontinuous clays are present. A full-scale demonstration, including 4% methane enhancement as a bioremediation nutrient in the injection well, was conducted during FY92, with results to be available in FY93. Better underground transport modeling and bioremediation modeling are needed. The technology was also used successfully in the DOE VOCs in the Non-Arid Soils Integrated Demonstration in Savannah River, South Carolina. Testing of directional boring for monitoring equipment installation was performed in an actual contamination zone during the summer of 1992.

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Cost:

Estimated costs are about \$60 to \$250 per meter (\$20 to \$75 per foot) for hydraulic bi-directional thrust drilling. Sonic drilling can be as much as \$330 per meter (\$100 per foot).

References:

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DOE, 1991. Horizontal Hybrid Directional Boring, FY92 Technical Task Plan, TTP Reference No.: AL-ZU23-J2.

DOE, 1991. SRS Integrated Demonstration: Directional Drilling, FY92 Technical Task Plan, TTP Reference No.: SR-1211-01.

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DOE, 1994. Technology Catalogue, First Edition. February.

DOE, 1993. Technology Name: Slant-Angle Sonic Drilling, Technology Information Profile (Rev. 2), DOE ProTech Database, TTP Reference No.: AL2310-05.

EPA, 1994. Manual: Alternative Methods for Fluid Delivery and Recovery, Prepared by: Murdoch, L., and Wilson, D.D. EPA/625/R-94/003.

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Kaback, D., and Oakley, D. 1996. "Horizontal Environmental Wells in the United States: A Catalogue," Colorado Center for Environmental Management (CCEM).

Wilson, D.D., and Kaback, D.S., 1993. "Industry Survey for Horizontal Wells", Westinghouse Savannah River Company (DOE), Aiken, SC. WSRC-TR-93-511.

Wilson, D.D., 1996. "Use of Horizontal Drilling In Environmental Remediation: A Horizontal Well Case Study", Water Well Journal, February.

Wilson, D.D., 1996, "Horizontal Well Development Made Easy", Water Well Journal, October.

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#### Site Information:

• DOE Demo Savannah River Site, SC

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**Points of Contact:** 



**Technology Specific Web Sites:** 



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# **Vendor Information:**

A list of vendors offering In Situ Physical/Chemical Water Treatment is available from the Vendor Information System for Innovative Treatment Technologies (VISITT) developed by U.S. Environmental Protection Agency (EPA).



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Health and Safety:



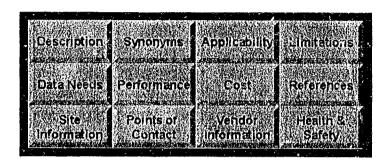
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# 4.58 Ground Water Pumping

(Ground Water Containment Remediation Technology)

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Technology	Description	
Ground Water, Surface Water, and Leachate		
3.13 Containment		
4.58 Ground Water Pumping	Ground water pumping is a component of many pump-and-Treat processes, which are some of the most commonly used ground water remediation technologies at contaminated sites.	

## Description:

Figure 4-58:

Typical Ground Water Pumping System Possible objectives of ground water pumping include removal of dissolved contaminants from the subsurface, and containment of contaminated ground water to prevent migration.

The first step of any remediation project consists of defining the remedial action objectives to be accomplished at the site. This involves gathering enough background site information and field data to make assessments of remedial requirements and possible cleanup levels. The first determination is whether cleanup or containment will be the most appropriate remedial action. If cleanup is chosen, the level of cleanup must be determined. If containment is chosen, ground water pumping is used as a hydraulic barrier to prevent off-site migration of contaminant plumes.

The next component consists of the design and implementation of the ground water pumping system based on data evaluated in setting the goals and objectives. The criteria for well design, pumping system, and treatment are dependent on the physical site characteristics and contaminant type. Actual treatment may include the design of a train of processes such as gravity segregation, air strippers, carbon systems tailored to remove specific contaminants.

Another component of any ground water extraction system is a ground water monitoring program to verify its effectiveness. Monitoring the remedial with wells and piezometers allows the operator to make iterative adjustments to the system in response to changes in subsurface conditions caused by the remediation.

The final component is determining the termination requirements. Termination requirements are based on the cleanup objectives defined in the initial stage of the remedial process. The termination criteria are also dependent on the specific site aspects revealed during remedial operations.

Although pumping for containment implies no treatment the following treatments usually follow pumping in pump and treat systems. These are briefly described below and in detail in technology profiles 4.47 through 4.55:

# 4.47 Bioreators:

Contaminants in extracted ground water are put into contact with microorganisms in attached or suspended growth biological reactors. In suspended systems, such as activated sludge, contaminated ground water is circulated in an aeration basin. In attached systems, such as rotating biological contractors and trickling filters, microorganisms are established on an inert support matrix.

## 4.48 Constructed wetlands:

The constructed wetlands-based treatment technology uses natural geochemical and biological processes inherent in an artificial wetland ecosystem to accumulate and remove metals and other contaminants from influent waters.

## 4.49 Adsorption/Absorption:

In liquid adsorption, solutes concentrate at the surface of a sorbent, thereby reducing their concentration in the bulk liquid phase. The most common adsorbent is granulated activated carbon (GAC) (see Technology Profile No. 4.51). Other natural and synthetic adsorbents include: forage sponge, lignin adsorption, sorption clays, and synthetic resins.

## 4.50 Air Stripping:

Volatile organics are partitioned from ground water by increasing the surface area of the contaminated water exposed to air. Aeration methods include packed towers, diffused aeration, tray aeration, and spray aeration.

# 4.51 Granulated Activated Carbon (GAC)/Liquid Phase Carbon Adsorption:

Ground water is pumped through a series of canisters or columns containing activated carbon to which dissolved organic contaminants adsorb. Periodic replacement or regeneration of saturated carbon is required.

4.52 Ion Exchange:

Ion exchange removes ions from the aqueous phase by the exchange of cations or anions between the contaminants and the exchange medium. Ion exchange materials may consist of resins made from synthetic organic materials that contain ionic functional groups to which exchangeable ions are attached. They also may be inorganic and natural polymeric materials. After the resin capacity has been exhausted, resins can be regenerated for re-use.

4.53 Precipitation/Coagulation/ Flocculation:

This process transforms dissolved contaminants into an insoluble solid, facilitating the contaminant's subsequent removal from the liquid phase by sedimentation or filtration. The process usually uses pH adjustment, addition of a chemical precipitant, and flocculation.

4.54 Separation:

Separation processes seek to detach contaminants from their medium (i.e., ground water and/or binding material that contain them). Ex situ separation of waste stream can be performed by many processes: (1) distillation, (2) filtration/ultrafiltration/microfiltration, (3) freeze crystallization, (4) membrane pervaporation and (5) reverse osmosis.

4.55 Sprinkler Irrigation:

Wastewater is distributed over the top of the filter bed through which wastewater is trickled. The organic contaminants in wastewater are degraded by the microorganisms attached to the filter medium.

# > Surfactant Enhanced Recovery

The application of surfactants micelles or steam to the ground water can facilitate the ground water pumping process by increasing the mobility and solubility of the contaminants sorbed to the soil matrix. This material can also facilitate the entrainment of hydrophobic contaminants to allow removal and assures that multi-phase contaminants can be effectively removed. Thus it can increase the contaminant mass removal per pore volume of ground water flushing through the contaminated zone.

The implementation of surfactant-enhanced recovery requires the injection of surfactants into a contaminated aquifer. Typical systems utilize a pump to extract ground water at some distance away from the injection point. The extracted ground water is treated ex situ to separate the injected surfactants from the contaminants and ground water. In order to be cost-effective, the design of the surfactant-enhanced recovery system is critical. Once the surfactants have separated from the ground water they can be re-injected into the subsurface. Contaminants must be separated from the ground water and treated prior to discharge of the extracted ground water.

# ➤ Drawdown Pumping

Pump drawdown nonaqueous-phase liquid (NAPL) recovery systems are designed to pump NAPL and ground water from recovery wells or trenches. Pumping removes water and lowers the water table near the extraction area to create a cone of depression. The cone of depression in the vicinity of the extraction well produces a gravity head that pushes flow of NAPL toward the well and increases the thickness of the NAPL layer in the well. Each foot of ground water depression provides a driving head equivalent to a pressure difference of 0.45 psi. In most cases, the production of a cone of depression will increase NAPL recovery rates.

Pumping may be accomplished with one or two pumps. In the single-pump configuration, one pump withdraws both water and NAPL. The dual-pump configuration uses one pump located below the water table to remove water and a second located in the NAPL layer to recover NAPL. A single-pump system reduces capital and operating costs and allows simpler control systems and operation, but produces a stream of mixed water and NAPL that must then be separated.

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Synonyms:

Pump and treat.

**DSERTS Code:** Q1 (Waste Removal - Liquids)

#### Dack to Tup

Applicability:

The first step in determining whether ground water pumping is an appropriate remedial technology is to conduct a site characterization investigation. Site characteristics, such as hydraulic conductivity, will determine the range of remedial options possible. Chemical properties of the site and plume need to be determined to characterize transport of the contaminant and evaluate the feasibility of ground water pumping. To determine if ground water pumping is appropriate for a site, one needs to know the history of the contamination event, the properties of the subsurface, and the biological and chemical contaminant characteristics. Identifying the chemical and physical site characteristics, locating the ground water contaminant plume in three dimensions, and determining aquifer and soil properties are necessary in designing an effective ground water pumping strategy.

Surfactant-enhanced recovery are most applicable for contaminated sites with enhanced light, nonaqueous-phase liquids (LNAPLs) and dense, nonaqueous-phase liquids (DNAPLs).

Drawdown pumping is effective for NAPL recovery when the aquifer has moderate to high hydraulic conductivity and a thick layer of low-viscosity NAPL. An aquifer with high hydraulic conductivity gives less flow resistance of NAPL into the well. A thick layer of NAPL allows the pumping system to collect a high proportion of NAPL in relation to the amount of ground water. For best operation, the NAPL thickness should be sufficient to completely cover the pump suction port.

Drawdown pumping is a commercially available technology that can be easily implemented with conventional pumps in wells or trenches. System installation costs are low to moderate, but the cost per amount of NAPL recovered varies greatly.

#### Dack to Ton

#### Limitations:

The following factors may limit the applicability and effectiveness of ground water pumping as part of the remedial process:

- The potentially long time necessary to achieve the remediation goal
- System designs fail to contain the contaminant as predicted, allowing the plume to migrate and failure of the pumping equipment.
- Residual saturation of the contaminant in the soil pores cannot be removed by ground water pumping. Con aminants tend to be sorbed in the soil matrix. Ground water pumping is not applicable to contaminants with high residual saturation, contaminants with high sorption capabilities, and homogeneous aquifers with hydraulic conductivity less than 10<sup>-5</sup> cm/sec.
- The cost of permitting procuring and operating treatment systems is high. Additional cost may also be attributed to the disposal of spend carbon and other treatment residuals and wastes.
- Biofouling of the extraction wells and associated treatment stream is a common problem which can severely affect system performance. The potential for this problem should be evaluated prior to the installation.

The following factors may limit the applicability and effectiveness of surfactant-enhanced recovery:

- Subsurface heterogeneities, as with most ground water remediation technologies, present challenges to the successful implementation of surfactant-enhanced recovery
- Potential toxic effects of residual surfactants in the subsurface
- Off-site migration of contaminants due to the increase solubility achieved with surfactant injection Obtaining regulatory approval to inject surfactants into an aquifer.

The following factors may limit the applicability and effectiveness of drawdown pumping:

• Drawdown pumping generally produces large volumes of water

in the process of recovering the free product.

• The production of a cone depression in the water table can smear the free product or trap the fuel in the saturated zone when the water table returns to its original level.

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Data Needs:

Collecting as much background site data as possible, obtaining accurate information on the type of contaminants present, and determining the hydrogeological nature of the site are essential. Contaminant information needed consists of: 1) source characterization, including the volume released, the area infiltrated, and duration of release; 2) concentration distribution of contaminants and naturally occurring chemicals in the ground water and soil; and 3) processes affecting plume development, such as chemical and biological reactions influencing contaminant mobility. Hydrogeologic information include determining the size of the contaminated aquifer, depth to water table, hydraulic conductivity of the surrounding aquifer material, impermeable units and confining layers, ground water flow direction and velocity, recharge and discharge areas, seasonal variations of ground water conditions, and local ground water use. Methods for determining aquifer properties include a slug test, pump test, and a borehole flowmeter test. The pump test consists of pumping one well and measuring the water level response of surrounding wells. A slug test measures the rate at which water level in one well returns to its initial state after inducing a rapid water level change by introducing or with drawing a volume of water. The borehole flowmeter test measures flow direction and rate in a borehole.

**Performance Data:** 

DOE has developed and tested many pump and treat technologies for hazardous waste removal over the past twenty years. Performance data on some of the most recent DOE sites can be in the Technology Application Analyses and site information below.

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Cost:

Cost data varies from site to site for ground water pump and treat technology. Recent cost data on a few of DOE's sites can be found below in the links to the Technology Application Analyses and in the site information.

Additional cost information can be found in the Hazardous, Toxic, and Radioactive Wastes (HTRW) Historical Cost Analysis System (HCAS) developed by Environmental Historical Cost Committee of Interagency Cost Estimation Group.

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- Pump and Treat System at Commencement Bay, South Tacoma Channel (Well 12A), Phase 2, Tacoma, Washington
- Recovery of Free Petroleum Product at Fort Drum, Fuel Dispensing Area 1595, Watertown, New York
- Pump and Treat of Contaminated Groundwater at IRP Site 4, Langley Air Force Base, Virginia
- Pump and Treat of Contaminated Groundwater at Operable Unit B/C, McClellan Air Force Base, California
- Pump and Treat of Contaminated Groundwater at the Twin Cities Army Ammunition Plant, New Brighton, Minnesota
- Pump and Treat of Contaminated Groundwater at the U.S. Department of Energy's Kansas City Plant, Kansas City, Missouri
- Pump and Treat of Contaminated Groundwater at the U.S.
   Department of Energy's Savannah River Site, A/M Area, Aiken, South Carolina

Federal Remediation Technologies Roundtable, 1998. Remediation Case Studies: Groundwater Pump and Treat (Chlorinated Solvents), EPA/542/R-98/013

- Pump and Treat of Contaminated Groundwater at the Des Moines TCE Superfund Site, OU 1, Des Moines, Iowa
- Pump and Treat of Contaminated Groundwater at the Former Firestone Facility Superfund Site, Salinas, California
- Pump and Treat of Contaminated Groundwater at the JMT.
   Facility RCRA Site, Brockport, New York
- Pump and Treat of Contaminated Groundwater at the Keefe Environmental Services Superfund Site, Epping, New Hampshire
- Groundwater Pump and Treat and Soil Vapor Extraction at DOE's Lawrence Livermore National Laboratory, Site 300,

GSA OU, Livermore, California

 Pump and Treat of Contaminated Groundwater at the Mystery Bridge at Hwy 20 Superfund Site, Dow/DSI Facility, Evansville, Wyoming

Groundwater Containment at Site LF-12, Offutt AFB, Nebraska

 Pump and Treat of Contaminated Groundwater at the Old Mill Superfund Site, Rock Creek, Ohio

 Pump and Treat of Contaminated Groundwater at the SCRDI Dixiana Superfund Site, Cayce, South Carolina

- Groundwater Containment at Site OT-16B, Shaw AFB, South Carolina
- Groundwater Containment at Site SD-29 and ST-30, Shaw AFB, South Carolina
- Pump and Treat of Contaminated Groundwater at the Sol Lynn/Industrial Transformers Superfund Site, Houston, Texas

 Pump and Treat of Contaminated Groundwater at the Solid State Circuits Superfund Site, Republic, Missouri

 Pump and Treat of Contaminated Groundwater with Containment Wall at the Solvent Recovery Services of New England, Inc. Superfund Site, Southington, Connecticut

Federal Remediation Technologies Roundtable, 1998. Remediation Case Studies: Groundwater Pump and Treat (Nonchlorinated Solvents), EPA/542/R-98/014

 Pump and Treat of Contaminated Groundwater at the Baird and McGuire Superfund Site, Holbrook, Massachusetts

 Pump and Treat of Contaminated Groundwater at the City Industries Superfund Site, Orlando, Florida

 Pump and Treat of Contaminated Groundwater at the King of Prussia Technical Corporation Superfund Site, Winslow Township, New Jersey

• Pump and Treat of Contaminated Groundwater at the LaSalle Electrical Superfund Site, LaSalle, Illinois

 Pump and Treat of Contaminated Groundwater at the Mid-South Wood Products Superfund Site, Mena, Arkansas

 Pump and Treat of Contaminated Groundwater at the Odessa Chromium J Superfund Site, OU 2, Odessa, Texas

 Pump and Treat of Contaminated Groundwater at the Odessa Chromium IIS Superfund Site, OU 2, Odessa, Texas

• Groundwater Containment at Site FT-01, Pope AFB, North Carolina

- Groundwater Containment at Site SS-07, Pope AFB, North Carolina
- Pump and Treat and Containment of Contaminated Groundwater at the Sylvester/Gilson Road Superfund Site, Nashua, New Hampshire

 Pump and Treat of Contaminated Groundwater at the United Chrome Superfund Site, Corvallis, Oregon

 Pump and Treat of Contaminated Groundwater at the U.S. Aviex Superfund Site, Niles, Michigan

 Pump and Treat of Contaminated Groundwater at the Western Processing Superfund Site, Kent, Washington Federal Remediation Technologies Roundtable, 1998. Remediation Case Studies: Innovative Groundwater Treatment Technologies, EPA/542/R-98/015.

 Pump and Treat and Permeable Reactive Barrier to Treat Contaminated Groundwater at the Former Intersil, Inc. Site, Sunnyvale, California

Pump and Treat and In Situ Bioremediation of Contaminated
 Groundwater at the French Ltd. Superfund Site, Crosby, Texas

 Pump and Treat and Air Sparging of Contaminated Groundwater at the Gold Coast Superfund Site, Miami, Florida

 Pump and Treat and In Situ Bioremediation of Contaminated Groundwater at the Libby Groundwater Superfund Site, Libby, Montana

 Pump and Treat, In Situ Bioremediation, and In Situ Air Sparging of Contaminated Groundwater at Site A, Long Island, New York

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#### **Site Information:**

- Lawerence Livermore Nat'l Lab (LLNL), Livermore, CA
- McClellan Air Force Base, Operable Unit (OU) D, CA
- FAA Technical Center, NJ
- Former Intersil, Inc. Site, Sunnyvale, CA.
- Amoco Petroleum Pipeline, Constantine, MI
- Baird and McGuire Superfund Site, Holbrook, MA
- City Industries Superfund Site, Orlando, FL
- Commencement Bay, South Tacoma Channel (Well 12A), Phase 2, Tacoma, WA
- Des Moines TCE Superfund Site, OU 1, Des Moines, IO
- Former Firestone Facility Superfund Site, Salinas, CA
- Fort Drum, Fuel Dispensing Area 1595, Watertown, NY
- French Ltd. Superfund Site, Crosby, TX
- Gold Coast Superfund Site, Miami, FL
- JMT Facility RCRA Site, Brockport, NY
- Keefe Environmental Services Superfund Site, Epping, NH
- King of Prussia Technical Corporation Superfund Site, Winslow Township, NJ
- Langley Air Force Base, IRP Site 4, Langley, VA
- LaSalle Electrical Superfund Site, LaSalle, IL
- DOE's Lawrence Livermore National Laboratory, Site 300, GSA OU, Livermore, CA
- Libby Groundwater Superfund Site, Libby, MT
- McClellan Air Force Base, Operable Unit (OU) B/C,CA
- · Mid-South Wood Products Superfund Site, Mena, AR
- Mystery Bridge at Hwy 20 Superfund Site, Dow/DSI Facility, Evansville, WY
- Odessa Chromium I Superfund Site, OU 2, Odessa, TX
- Odessa Chromium II.3 Superfund Site, OU 2, Odessa, TX
- Site LF-12, Offutt AFB, NB
- Old Mill Superfund Site, Rock Creek, OH

- Site FT-01, Pope AFB, NC
- Site SS-07, Pope AFB, NC
- SCRDI Dixiana Superfund Site, Cayce, SC
- Site OT-16B, Shaw AFB, SC
- Site SD-29 and ST-30, Shaw AFB, SC
- Site A, Long Island, NY
- Sol Lynn/Industrial Transformers Superfund Site, Houston, TX
- Solid State Circuits Superfund Site, Republic, MO
- Solvent Recovery Services of New England, Inc. Superfund Site, Southington, CT
- Sylvester/Gilson Road Superfund Site, Nashua, NH
- Twin City Army Ammunition Plant, New Brighton, MN
- United Chrome Superfund Site, Corvallis, OR
- U.S. Aviex Superfund Site, Niles, MI
- U.S. DOE's Kansas City Plant, Kansas City, MO
- U.S. DOE's Savannah River Site A/M Area, Aiken, SC
- Western Processing Superfund Site, Kent, WA
- Additional site information on the FRTR web site

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**Points of Contact:** 



**Technology Specific Web Sites:** 



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#### Vendor Information:

A list of vendors offering Water Containment Treatment is available from the Vendor Information System for Innovative Treatment Technologies (VISITT) developed by U.S. Environmental Protection Agency (EPA).



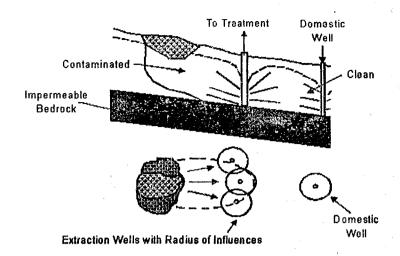
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# 4.59 Slurry Walls

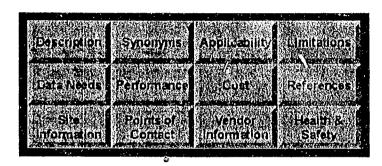
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Technology	<b>D</b> escription	
Ground Water, Surface Water, and Leachate		
3.13 Containment		
	These subsurface barriers consist of vertically excavated trenches filled with slurry. The slurry, usually a mixture of bentonite and water, hydraulically shores the trench to prevent collapse and retards ground water flow.	

**Description:** 

Figure 4-59: Typical Keyed-In Slurry Wall (Cross-Section)

Slurry walls are used to contain contaminated ground water, divert contaminated ground water from the drinking water intake, divert uncontaminated ground water flow, and/or provide a barrier for the ground water treatment system.

These subsurface barriers consist of a vertically excavated trench that is filled with a slurry. The slurry hydraulically shores the trench to prevent collapse and forms a filter cake to reduce ground water flow. Slurry walls often are used where the waste mass is too large for treatment and where soluble and mobile constituents pose an imminent threat to a source of drinking water.

Slurry walls are a full-scale technology that have been used for decades as long-term solutions for controlling seepage. They are often used in conjunction with capping. The technology has demonstrated its effectiveness in containing greater than 95% of the uncontaminated ground water; however, in contaminated ground water applications, specific contaminant types may degrade the slurry wall components and reduce the long-term effectiveness.

Most slurry walls are constructed of a soil, bentonite, and water

mixture. The bentonite slurry is used primarily for wall stabilization during trench excavation. A soil-bentonite backfill material is then placed into the trench (displacing the slurry) to create the cutoff wall. Walls of this composition provide a barrier with low permeability and chemical resistance at low cost. Other wall compositions, such as cement/bentonite, pozzolan/bentonite, attapulgite, organically modified bentonite, or slurry/geomembrane composite, may be used if greater structural strength is required or if chemical incompatibilities between bentonite and site contaminants exist.

Slurry walls are typically placed at depths up to 30 meters (100 feet) and are generally 0.6 to 1.2 meters (2 to 4 feet) in thickness. Installation depths over 30 m (100 ft) are implementable using clam shell bucket excavation, but the cost perunit area of wall increases by about a factor of three. The most effective application of the slurry wall for site remediation or pollution control is to base (or key) the slurry wall 0.6 to 0.9 meters (2 to 3 feet) into a low permeability layer such as clay or bedrock, as shown in the preceding figure. This "keying-in" provides for an effective foundation with minimum leakage potential. An alternate configuration for slurry wall installation is a "hanging" wall in which the wall projects into the ground water table to block the movement of lower density or floating contaminants such as oils, fuels, or gases. Hanging walls are used less frequently than keyed-in walls.

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Synonyms:

Vertical cutoff walls; Hydrodynamic barriers; Slurry Trenches. **DSERTS Code:** 12 (Slurry Walls/Underground Barriers)

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Applicability:

Slurry walls contain the ground water itself, thus treating no particular target group of contaminants. They are used to contain contaminated ground water, divert contaminated ground water from drinking water intake, divert uncontaminated ground water flow, and/or provide a barrier for the ground water treatment system.

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Limitations:

Factors that may limit the applicability and effectiveness of the process include:

- Most of the approaches involve a large amount of heavy construction.
- The technology only contains contaminants within a specific area.
- Soil-bentonite backfills are not able to withstand attack by strong acids, bases, salt solutions, and some organic chemicals. Other slurry mixtures can be developed to resist specific chemicals.
- There is the potential for the slurry walls to degrade or deteriorate over time.
- Use of this technology does not guarantee that further remediation in the future may not be necessary.

# Data Needs:

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A detailed discussion of these data elements is provided in <u>Subsection 2.2.2</u>. (Data Requirements for Ground Water, Surface Water, and Leachate).

The following factors, at a minimum, must be assessed prior to designing effective soil-bentonite slurry walls: maximum allowable permeability, anticipated hydraulic gradients, required wall strength, availability and grade of bentonite to be used, boundaries of contamination, compatibility of wastes and contaminants in contact with slurry wall materials, characteristics (i.e., depth, permeability, and continuity) of substrate into which the wall is to be keyed, characteristics of backfill material (e.g., fines content), and site terrain and physical layout.

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#### Performance Data: \*

Slurry walls have been used for decades, so the equipment and methodology are readily available and well known; however, the process of designing the proper mix of wall materials to contain specific contaminants is less well developed. Excavation and backfilling of the trench is critical and requires experienced contractors.

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#### Cost:

Costs likely to be incurred in the design and installation of a standard soil-bentonite wall in soft to medium soil range from \$540 to \$750 per square meter (\$5 to \$7 per square foot) (1991 dollars). These costs do not include variable costs required for chemical analyses, feasibility, or compatibility testing. Testing costs depend heavily on site-specific factors.

Factors that have the most significant impact on the final cost of soilbentonite slurry wall installation include:

- Type, activity, and distribution of contaminants.
- Depth, length, and width of wall.
- Geological and hydrological characteristics.
- Distance from source of materials and equipment.
- Requirements for wall protection and maintenance.
- Type of slurry and backfill used.
- Other site-specific requirements as identified in the initial site assessment (e.g., presence of contaminants or debris).
- Planning, permitting, regulatory interaction, and site restoration.

Additional cost information can be found in the Hazardous, Toxic, and Radioactive Wastes (HTRW) Historical Cost Analysis System (HCAS) developed by Environmental Historical Cost Committee of Interagency Cost Estimation Group.

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## **Site Information:**

- Hazardous Waste Landfill
- Sanitary Landfill
- Coal Tar Disposal Pond
- Sylvester/Gilson Road Superfund Site, Nashua, NH

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# **Points of Contact:**



# **Technology Specific Web Site:**



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## **Vendor Information:**

A list of vendors offering In Situ Physical/Chemical Water Treatment is available from the Vendor Information System for Innovative Treatment Technologies (VISITT) developed by U.S. Environmental Protection Agency (EPA).



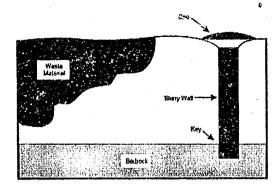
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## Health and Safety:



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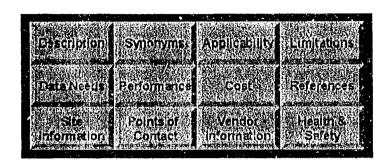
# 4.30 Landfill Cap

(Soil Containment Remediation Technology)

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Technology	Description	
Soil, Sediment, and Sludge		
3.7 Containment		
4.30 Landfill Cap	Landfill caps are used for contaminant source control.	

Description:

Landfill caps can be used to:

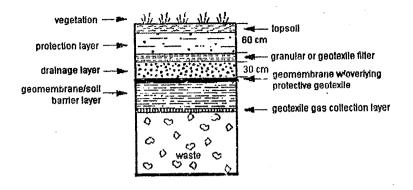
Figure 4-30: Typical

Figure 4-30: Typical RCRA Subtitle C Landfill Cap System

- Caps can minimize exposure on the surface of the waste facility.
- Prevent vertical infiltration of water into wastes that would create contaminated leachate.
- Contain waste while treatment is being applied.
- Control gas emissions from underlying waste.
- Create a land surface that can support vegetation and/or be used for other purposes.

Landfill Capping is the most common form of remediation because it is generally less expensive than other technologies and effectively manages the human and ecological risks associated with a remediation site.

The design of landfill caps is site specific and depends on the intended functions of the system. Landfill Caps can range from a one-layer system of vegetated soil to a complex multi-layer system of soils and geosynthetics. In general, less complex systems are required in dry climates and more complex systems are required in wet climates. The material used in the construction of landfill caps include low-permeability and high-permeability soils and low-permeability geosynthetic products. The low-permeability materials divert water



and prevent its passage into the waste. The high permeability materials carry water away that percolates into the cap. Other materials may be used to increase slope stability.

The most critical components of a landfill cap are the barrier layer and the drainage layer. The barrier layer can be low-permeability soil (clay) and/or geosynthetic clay liners (GCLs). A flexible geomembrane liner is placed on top of the barrier layer. Geomembranes are usually supplied in large rolls and are available in several thickness (20 to 140 mil), widths (15 to 100 ft), and lengths (180 to 840 ft). The candidate list of polymers commonly used is lengthy, which includes polyvinyl chloride (PVC), polyethylenes of various densities, reinforced chlorosulfonated polyethylene (CSPE-R), polypropylene, ethylene interpolymer alloy (EIA), and many newcomers. Soils used as barrier materials generally are clays that are compacted to a hydraulic conductivity no greater than 1 x 10<sup>-6</sup> cm/sec. Compacted soil barriers are generally installed in 6-inch minimum lifts to achieve a thickness of 2 feet or more. A composite barrier uses both soil and a geomembrane, taking advantage of the properties of each. The geomembrane is essentially impermeable, but, if it develops a leak, the soil component prevents significant leakage into the underlying waste.

For facilities on top of putrescible wastes, the collection and control of methane and carbon dioxide, potent greenhouse gases, must be part of facility design and operation.

# > Asphalt/Concrete Cap

The most effective single-layer caps are composed of concrete or bituminous asphalt. It is used to form a surface barrier between landfill and the environment. An asphalt concrete cap would reduce leaching through the landfill into an adjacent aquifer.

# > RCRA Subtitle C Cap

The RCRA C multilayered landfill cap is a baseline design that is suggested for use in RCRA hazardous waste applications. These caps generally consist of an upper vegetative (topsoil) layer, a drainage layer, and a low permeability layer which consists of a synthetic liner over 2 feet of compacted clay. The compacted clay liners are effective if they retain a certain moisture content but are susceptible to cracking if the clay material is desiccated. As a result alternate cap designs are usually considered for arid environments.

# > RCRA Subtitle D Cap

RCRA Subtitle D requirements are for non-hazardous waste landfills. The design of a landfill cover for a RCRA Subtitle D facility is generally a function of the bottom liner system or natural subsoils present. The cover must meet the following specifications:

- the material must have a permeability no greater than 1 x 10-5 cm/s, or equivalent permeability of any bottom liner or natural subsoils present, whichever is less.
- The infiltration layer must contain at least 45 cm of earthen material.
- The erosion control layer must be at least 15 cm of earthen material capable of sustaining native plant growth.

Alternative design can be considered, but must be of equivalent performance as the specifications outlined above. All covers should be designed to prevent the "bathtub" effect. The bathtub effect occurs when a more permeable cover is placed over a less permeable bottom liner or natural subsoil. The landfill then fills up like a bathtub.

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Synonyms:

Cap; Landfill cover; Surface cover.

**DSERTS Codes:** 

I0 (Containment)
I1 (Capping)

#### . Back to Top:

Applicability:

Landfill Caps may be temporary or final. Temporary caps can be installed before final closure to minimize generation of leachate until a better remedy is selected. They are usually used to minimize infiltration when the underlying waste mass is undergoing settling. A more stable base will thus be provided for the final cover, reducing the cost of the post-closure maintenance. Landfill caps also may be applied to waste masses that are so large that other treatment is impractical. At mining sites for example, caps can be used to minimize the infiltration of water to contaminated tailings piles and to provide a suitable base for the establishment of vegetation. In conjunction with water diversion and detention structures, landfill caps may be designed to route surface water away from the waste area while minimizing erosion.

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Limitations:

Landfilling does not lessen toxicity, mobility, or volume of hazardous wastes, but does mitigate migration. Landfill caps are most effective where most of the underlying waste is above the water table. A cap, by itself, cannot prevent the horizontal flow of ground water through the waste, only the vertical entry of water into the waste. In many cases landfill caps are used in conjunction with vertical walls to minimize horizontal flow and migration. The effective life of landfill components (including cap) can be extended by long-term inspection and maintenance. Vegetation, which has a tendency for deep root penetration, must be eliminated from the cap area. In addition, precautions must be taken to assume that the integrity of the cap is not compromised by land use activities.

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Data Needs:

A detailed discussion of these data elements is provided in Subsection 2.2.1 (Data Requirements for Soil, Sediment, and Sludge). Many laboratory tests are needed to ensure that the materials being considered for each of the landfill cap components are suitable. Tests to determine the suitability of soil include grain size analysis, Atterberg limits, and compaction characteristics. Landfill instability can be solved by understanding interface friction properties between all material layers, natural or synthetic. The major engineering soil properties that must be defined are the shear strength and hydraulic conductivity. Shear strength may be determined with the unconfined compression test, direct shear test, or triaxial compression test. Hydraulic conductivity of soils may be measured in the laboratory by the constant head permeability test or the falling head permeability test. Field hydraulic conductivity tests on test pads are generally recommended prior to actual cover construction to ensure that the low-permeability requirements can actually be met under construction conditions.

Laboratory tests are also needed to ensure that geosynthetic materials will meet the cap requirements, For example, geosynthetics in caps may be subjected to tensile stresses caused by subsidence and by the gravitational tendency of a geomembrane or material adjacent to it to slide or be pulled down slopes.

Since facility performance is a function of quality construction more so than selection of materials, construction quality assurance of caps are critical. EPA has generated a technical guidance document on this subject. The technical guidance should be strictly followed during design and construction.

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Performance Data:

Previously installed caps are hard to monitor for performance. Monitoring well systems or infiltration monitoring systems can provide some information, but it is often not possible to determine whether the water or leachate originated as surface water or ground water. Performance can be monitored much more effectively by including pan lysimeter in future caps.

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Cost:

Landfill caps are generally the least expensive way to manage the human health and ecological risks effectively. Rough industry cost are \$175k/acre for RCRA Subtitle D, and \$225k/acre for RCRA Subtitle C.

Additional cost information can be found in the Hazardous, Toxic, and Radioactive Wastes (HTRW) Historical Cost Analysis System (HCAS) developed by Environmental Historical Cost Committee of Interagency Cost Estimation Group.

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References:

DOE, 1995. *Technology Catalogue*, *Second Edition*, Office of Environmental Management and Office of Technology Development, DOE/EM-0235.

EPA, 1987. Geosynthetic Design Guidance for Hazardous Waste Landfill Cells and Surface Impoundments, EPA/600/2-87/025.

EPA, 1988. Technology Screening Guide for Treatment of CERCLA Soil and Sludge, EPA/540/288/004.

EPA, 1989. Final Covers on Hazardous Waste Landfills and Surface Impoundments, Technical Guidance Document, Office of Solid Waste and Emergency Response, Washington, DC, EPA/530/SW-89/047.

EPA, 1991. Compilation of Information on Alternative Barriers for Liner and Cover Systems, EPA/600/2-91/002.

EPA, 1991. Inspection Techniques for the Fabrication of Geomembrane Field Seams, Technical Guidance Document, EPA/530/SW-91/051.

EPA, 1993. Construction Quality Control and Quality Assurance at Waste Containment Facilities, Technical Guidance Document, Office of Research and Development, RREL, Cincinnati, OH, EPA/600/R-93/182.

EPA, 1993. Engineering Bulletin: Landfill Covers, EPA/540/S-93/500.

EPA, 1995. RCRA Subtitle D (258) Seismic Design Guidance for Municipal Solid Waste Landfills, Office of Research and Development, NRMRL, Cincinnati, OH, EPA/600/R-95/051.

EPA, 1995. Report of 1995 Workshop on Geosynthetic Clay Liners, Office of Research and Development, NRMRL, Cincinnati, OH, EPA/600/R-96/149.

EPA, 1997. Best Management Practices (BMPs) for Soil Treatment Technologies: Suggested Operational Guidelines to Prevent Crossmedia Transfer of Contaminants During Clean-UP Activities, EPA OSWER, EPA/530/R-97/007.

EPA, 1997. Technology Alternatives for the Remediation of Soils Contaminated with As, Cd, Cr, IIg, and Pb, Engineering Bulletin, EPA540/R-97/008.

EPA, 1998. Evaluation of Subsurface Engineered Barriers at Waste Sites, Technology Report, EPA/542/R-98/005.

Federal Remediation Technologies Roundtable, 1998. Remediation

Case Studies: In Situ Soil Treatment Technologies (Soil Vapor Extraction, Thermal Processes), EPA/542/R-98/012

- Soil Vapor Extraction at the Seymour Recycling Corporation Superfund Site, Seymour, Indiana
- Federal Remediation Technologies Roundtable, 1998. Remediation
   Case Studies: Groundwater Pump and Treat (Nonchlorinated Solvents), EPA/542/R-98/014
  - Pump and Treat and Containment of Contaminated Groundwater at the Sylvester/Gilson Road Superfund Site, Nashua, New Hampshire

Federal Remediation Technologies Roundtable, 1998. Remediation Case Studies: Debris and Surface Cleaning Technologies, and Other Miscellaneous Technologies, EPA/542/R-98/017.

• Lawrence Livermore National Laboratory (LLNL) Site 300 - Pit 6 Landfill Operable Unit (OU), Livermore, CA.

#### Dack to Top

#### Site Information:

- · AFCEE action, Fairchild AFB, WA
- DOE Oak Ridge, TN facility
- Lawrence Livermore National Laboratory, Site 300, Coast Ranges, CA
- DOE Demo, Lee Acres landfill, Farmington, NM
- Lawrence Livermore National Laboratory, Site 300, Pit 6 Landfill OU, Livermore, CA
- Seymour Recycling Corporation Superfund Site, Seymour, IN
- Sylvester/Gilson Road Superfund Site, Nashua, NH

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#### **Points of Contact:**



#### **Technology Specific Web Sites:**



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#### Vendor Information:

A list of vendors offering Soil Containment Treatment is available from the Vendor Information System for Innovative Treatment Technologies (VISITT) developed by U.S. Environmental Protection Agency (EPA).



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Health and Safety:



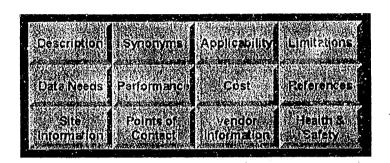
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# 4.31 Landfill Cap Enhancements

(Soil Containment Remediation Technology)

Synonym: .. Next Top Table of Previous Screen. Section List Section Page. Matrix Contents



Technology	Description			
Soil, Sediment, and Sludge				
3.7 Containment				
4.31 Landfill Cover Enhancements	The purpose of landfill cover enhancement is to reduce or eliminate contaminant migration (e.g. percolation). Water harvesting and vegetative cover are two ways for landfill cover enhancements. Water harvesting uses runoff enhancement to manage landfill site water balance. Vegetative cover reduces soil moisture via plant uptake and evapotranspiration.			

#### Description:

Figure 4-31: Typical

Water Harvesting Enhancement

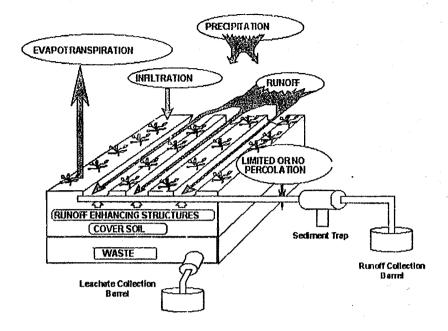
The precipitation to the landfill cover is balanced by the combination of following effects: run-off, cumulating in soil, evapotranspiration, and percolation. For a given amount of precipitation, in order to reduce or eliminate percolation, the effects of run-off and/or evapotranspiration need to be enhanced.

## ➤ Water Harvesting

Water harvesting uses runoff enhancement to manage landfill site water balance. This enhancement can be achieved by simply covering landfill cover surface with metal rain gutter placed parallel to the slope. The percentage of runoff increases when gutter coverage increases. However, too much coverage (> 40%) has little effect on runoff enhancement.

# > Vegetative Cover

Vegetative cover reduces soil moisture via plant uptake and evapotranspiration. Plant cover also limits soil erosion. Vegetative



cover is more stable because it emphasizes use of natural materials and configurations, which implies longevity.

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**Synonyms:** 

NA

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Applicability:

Landfill cover enhancement is applicable for traditional landfills, surface impoundment's, waste piles, sludges, and some mine tailings. It may prove to be less costly than a conventional barrier because it uses simple structure or local resources. It is simple in design, easy to install over an existing landfill cover, and easy to remove if other uses for the land emerge in the future.

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Limitations:

Factors that may limit the applicability and effectiveness of these processes include:

- Proper site evaluation is required.
- Plant coverage is seasonal.
- Too much gutter coverage (> 40%) has little effect on runoff enhancement.

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**Data Needs:** 

A detailed discussion of data elements is provided in <u>Subsection 2.2.1</u> (Data Requirements for Soil, Sediment, and Sludge).

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Performance Data:

Landfill cover enhancement is a fairly new technology that is still being tested and demonstrated.

Naval Facilities Engineering Services Center initiated a demonstration project at Marine Corp Base Hawaii to evaluate infiltration control cover design. Metal rain gutters were placed on the ground surface and parallel to the slope. Remaining surface was seeded with 6 native grasses and shrubs. After 9 months of operation, it was found that plots with runoff enhancement had 2 to 5 times more runoff, and 2 to 3 times less percolation.

#### Dack to Top.

Cost:

The simple configuration of landfill cover enhancement should result better containment with little increase in costs. These costs are currently determined on a case-by-case basis because of construction material availability and design requirements at various site locations.

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#### References:

Dwyer, S., 1997. *Alternative Landfill Cover Demonstration*. Sandia National Laboratories, U.S. DOE Environmental Management, Office of Science and Technology.

EPA, 1993. Evaluation of Subsurface Engineered Barriers at Waste Sites, Technology Report, EPA/542/R-98/005.

Finley, R., 1997. *Demonstrate Capillary Barrier Design Tools*. Sandia National Laboratories, U.S. DOE Environmental Management, Office of Science and Technology.

Hakonson, T.E., L. Karr, and B. Harre, 1997. A Water Balance Study of Infiltration Control Landfill Cover Design at Marine Corp Base Hawaii, Technical paper, Naval Facilities Engineering Services Center, Port Hueneme, CA.

Murphy B., 1995. Demonstration project may lead to more effective hazardous waste landfill covers. Sandia National Laboratories, U.S. DOE Environmental Management, Office of Science and Technology.

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#### Site Information:

- DOE Demonstration, Sandia National Laboratories
- MCAS Kaneohe Bay

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Points of Contact:



**Technology Specific Web Site:** 



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#### **Vendor Information:**

A list of vendors offering Soil Containment Treatment is available from the Vendor Information System for Innovative Treatment Technologies (VISITT) developed by U.S. Environmental Protection Agency (EPA).



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Health and Safety:



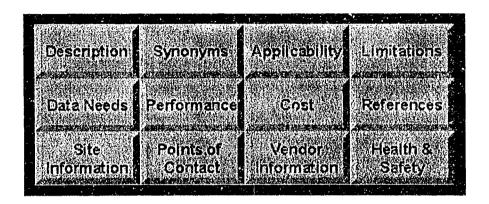
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# 4.32 Excavation, Retrieval, and Off-Site Disposal

(Other Soil Remediation Technology)

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Technology	Description			
Soil, Sediment, and Sludge				
3.7 Other Treatment				
4.32 Excavation, Retrieval, and Off-Site Disposal	Contaminated material is removed and transported to permitted off-site treatment and disposal facilities. Pretreatment may be required.			

#### Description:

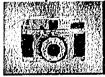


Figure 4-32: Typical Contaminated Soil Excavation Diagram

Contaminated material is removed and transported to permitted off-site treatment and/or disposal facilities. Some pretreatment of the contaminated media usually is required in order to meet land disposal restrictions.

Confined disposal facilities (CDFs) are engineered structure enclosed by dikes and designed to retain dredged materials. A CDF may have a large cell for material disposal, and adjoining cells for retention and decantation of turbid, supernatant water. A variety of linings have been used to prevent seepage through the dike walls. The most effective are clay or bentonite-cement slurries, but sand, soil, and sediment linings have also been used.

Location and design are two important CDF consideration. Terms to consider in the location of a CDF are the physical aspects (size, proximity to a navigable waterway), the design/construction (geology/hydrology), and the environmental (current use of the area, environmental value, and environmental effects). The primary goal of a CDF design is minimization of contaminant loss. Caps are the most effective way to minimize contaminant loss from CDFs, but selection of proper liner material is also an important control on CDFs. Finally, CDFs require continuous monitoring to ensure

structural integrity.

Operation and maintenance duration lasts as long as the life of the facility.

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Synonyms:

**DSERTS Codes:** 

E0 (Removal)

E1 (Waste Removal-Soils)
R1 (Waste Removal-Sludges)

S1 (Waste Removal-Non-soil Solids)

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Applicability:

Excavation and off-site disposal is applicable to the complete range of contaminant groups with no particular target group. Excavation and off-site by relocating the waste to a different (and presumably safer) site.

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Limitations:

Factors that may limit the applicability and effectiveness of the process include:

- Generation of fugitive emissions may be a problem during operations.
- The distance from the contaminated site to the nearest disposal facility with the required permit(s) will affect cost.
- Depth and composition of the media requiring excavation must be considered.
- Transportation of the soil through populated areas may affect community acceptability.
- Disposal options for certain waste (e.g., mixed waste or transuranic waste) may be limited. There is currently only one licensed disposal facility for radioactive and mixed waste in the United States.
- Contaminants can potentially migrate from CDF from several pathways, including effluent discharge to surface water, rainfall surface runoff, leachate into ground water, volatilization to the atmosphere, and dike uptake.
- CDFs can develop odor problems as well as mosquito and insect problems without proper design and maintenance.

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Data Needs:

A detailed discussion of these data elements is provided in Subsection 2.2.1 (Data Requirements for Soil, Sediment, and Sludge).

The type of contaminant and its concentration will impact off-site disposal requirements. Soil characterization as dictated by land disposal restrictions (LDRs) are required. Most hazardous wastes must be treated to meet either RCRA or non-RCRA treatment standards prior to land disposal. Radioactive wastes would have to meet disposal facility waste form requirements based on waste classification.

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Performance Data:

Excavation and off-site disposal is a well proven and readily implementable technology. Prior to 1984, excavation and off-site disposal was the most

common method for cleaning up hazardous waste sites. Excavation is the initial component in all ex situ treatments.

The rate of excavation depends on a number of factors, including the number of loaders and trucks operating. The excavation of 18,200 metric tons (20,000 tons) of contaminated soil would typically require about 2 months. Disposal of the contaminated media is dependent upon the availability of adequate containers to transport the hazardous waste to a permitted facility.

CERCLA includes a statutory preference for treatment of contaminants, and excavation and off-site disposal is now less acceptable than in the past. The disposal of hazardous wastes is governed by RCRA (40 CFR Parts 261-265), and the U.S. Department of Transportation (DOT) regulates the transport of hazardous materials (49 CFR Parts 172-179, 49 CFR Part 1387, and DOT-E 8876).

DOE has demonstrated a cryogenic retrieval of buried waste system, which uses liquid nitrogen (LN2) to freeze soil and buried waste to reduce the spread of contamination while the buried material is retrieved with a series of remotely operated tools. Other excavation/retrieval systems that DOE is currently developing include a remote excavation system, a hydraulic impact end effector, and a high pressure waterjet dislodging and conveyance end effector using confined sluicing.

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Cost:

Cost estimates for excavation and disposal range from \$300 to \$510 per metric ton (\$270 to \$460 per ton) depending on the nature of hazardous materials and methods of excavation. These estimates include excavation/removal, transportation, and disposal at a RCRA permitted facility. Additional cost of treatment at disposal facility may also be required. Excavation and off-site disposal is a relatively simple process, with proven procedures. It is a labor-intensive practice with little potential for further automation. Additional costs may include soil characterization and treatment to meet land ban requirements.

Additional cost information can be found in the Hazardous, Toxic, and Radioactive Wastes (HTRW) Historical Cost Analysis System (HCAS) developed by Environmental Historical Cost Committee of Interagency Cost Estimation Group.

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References:

Church, H.K., 1981. *Excavation Handbook*, McGraw Hill Book Co., New York, NY.

DOE, 1995. *Technology Catalogue*, *Second Edition*, Office of Environmental Management and Office of Technology Development, DOE/EM-0235.

EPA, 1991. Survey of Materials-Handling Technologies Used at Hazardous Waste Sites, EPA, ORD, Washington, DC, EPA/540/2-91/010.

EPA, 1992. McColl Superfund Site Demonstration of a Trial Excavation, EPA RREL, series include Technology Evaluation EPA/S40/R-92/015, PB92-226448; Applications Analysis, EPA/540/AR-92/015; and Technology Demonstration. Summary, EPA/540/SR/-92/015.

EPA, 1997. Best Management Practices (BMPs) for Soil Treatment Technologies: Suggested Operational Guidelines to Prevent Cross-media Transfer of Contaminants During Clean-UP Activities, EPA OSWER, EPA/530/R-97/007.

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#### Site Information:

- EPA Demo: Carter Industrial, MI; Shaver's Farm, GA; Hopkinsville, KY
- EPA Demo: IN, MI, OH, SD, VA, WI
- DOE Integrated Demo: (1,2) Chemical and Mixed Waste Landfills, Albuquerque, NM; (3) Mixed Waste Landfill at Kirkland AFB, NM
- DOE Integrated Demo: Fernald Environmental Project Cincinnati, OH

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#### **Points of Contact:**



## **Technology Specific Web Sites:**



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#### **Vendor Information:**

A list of vendors offering Soil Excavation, Retrieval, and Off-Site Disposal Treatment is available from the Vendor Information System for Innovative Treatment Technologies (VISITT) developed by U.S. Environmental Protection Agency (EPA).



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Health and Safety:



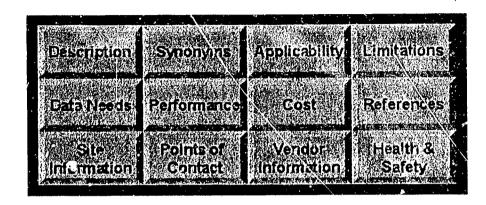
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# 4.36 Phytoremediation

(In Situ Ground Water Remediation Technology)

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Technology	Description		
Ground Water, Surface Water, and Leachate			
3.9 In Situ Biological Treatment			
4.36 Phytoremediation	Phytoremediation is a set of processes that uses plants to remove, transfer, stabilize and destroy organic/inorganic contamination in ground water, surface water, and leachate.		

### **Description:**



Figure 4-36 Typical In Situ Phytoremediation System

Phytoremediation is a set of processes that uses plants to clean contamination in ground water and surface water. The treatment of metals or other inorganic contamination has been discussed in Section 4.5 (Phytoremediation for Soil). There are several ways plants can be used for the phytoremediation. These mechanisms include enhanced rhizosphere biodegradation, hydraulic control, phyto-degradation and phyto-volatilization.

# > Enhanced Rhizosphere Biodegradation

Enhanced rhizosphere biodegradation takes place in the soil surrounding plant roots. Natural substances released by plant roots supply nutrients to microorganisms, which enhances their ability to biodegrade organic contaminants. Plant roots also loosen the soil and then die, leaving paths for transport of water and aeration. This process tends to pull water to the surface zone and dry the lower saturated zones.

# > Hydraulic Control

Depending on the type of trees, climate, and season, trees can act as organic

pumps when their roots reach down towards the water table and establish a dense root mass that takes up large quantities of water.

## > Phyto-degradation

Phyto-degradation is the metabolism of contaminants within plant tissues. Plants produce enzymes, such as dehalogenase and oxygenase, that help catalyze degradation. Investigations are proceeding to determine if both aromatic and chlorinated aliphatic compounds are amenable to phyto-degradation.

# > Phyto-volatilization

Phyto-volatilization occurs as plants take up water containing organic contaminants and release the contaminants into the air through their leaves. Plants can also break down organic contaminants and release breakdown products into air through leaves.

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Synonyms:

Vegetation-enhanced bioremediation.

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Applicability:

Phytoremediation can be used to clean up organic contaminants from surface water, ground water, leachate, and municipal and industrial wastewater.

Plants also produce enzymes, such as dehalogenase and oxygenase, which help catalyze degradation.

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#### Limitations:

There are a number of limitations to phytoremediation

- It is limited to shallow soils, streams, and ground water.
- High concentrations of hazardous materials can be toxic to plants.
- It involves the same mass transfer limitations as other biotreatments.
- Climatic or seasonal conditions may interfere or inhibit plant growth, slow remediation efforts, or increase the length of the treatment period.
- It can transfer contamination across media, e.g., from soil to air.
- It is not effective for strongly sorbed (e.g., PCBs) and weakly sorbed contaminants.
- Phytoremediation will likely require a large surface area of land for remediation.
- The toxicity and bioavailability of biodegradation products is not always known. Products may be mobilized into ground water or bioaccumulated in animals. More research is needed to determine the fate of various compounds in the plant metabolic cycle to ensure that plant droppings and products manufactured by plants do not contribute toxic or harmful chemicals into the food chain or increase risk exposure to the general public.

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Data Needs:

A detailed discussion of these data elements is provided in <u>Subsection 2.2.2</u>. (Data Requirements for Ground Water, Surface Water, and Leachate).

In addition, detailed information is needed to determine the kinds of soil used for phytoremediation projects. Water movement, reductive oxygen concentrations, root growth, and root structure all affect the growth of plants and should be considered when implementing phytoremediation.

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Performance Data:

The U.S. Air Force used poplar trees to contain a ground water TCE plume. TCE was shown to be degraded in the tissues of the poplar trees. The trees pumped a sufficient amount of water to produce a cone of depression limiting the spread of the TCE plume.

In Iowa, EPA demonstrated the usage of phytoremediation by planting poplar trees along a stream bank between a corn field and the stream. These trees acted as natural pumps to keep toxic herbicides, pesticides, and fertilizers out of the streams and ground water. After three years, while the nitrate concentration in ground water at the edge of the corn field was measured at 150 mg/L, the ground water among the poplar trees along the stream bank had nitrate concentration of only 3 mg/L.

USAEC is also leading the team of experts from EPA, Tennessee Valley Authority (TVA) and the Waterways Experimental Station (WES) to successfully demonstrate phytoremediation of explosive contaminated sites in Milan Army Ammunition Plant in Milan, TN.

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Cost:

Construction estimates for phytoremediation are \$200K/acre and \$20K/acre for operations and maintenance.

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References:

Boyajian, G. E. and Devedjian, D. L., 1997. "Phytoremediation: It Grows on You", Soil & Groundwater Cleanup, February/March, pp. 22-26.

California Base Closure Environmental Committee (CBCEC), 1994.

Treatment Technologies Applications Matrix for Base Closure Activities,
Revision 1, Technology Matching Process Action Team, November, 1994.

EPA, 1998. A Citizen's Guide to Phytoremediation, Technology Fact Sheet, EPA NCEPI, EPA/542/F-98/011.

EPA, 1996. A Citizen's Guide to Bioremediation, Technology Fact Sheet, EPA NCEPI, EPA/542/F-96/007.

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USAFC, 1997. Phytoremediation of Explosives in Groundwater Using Constructed Wetlands in Innovative Technology Demonstration, Evaluation and Transfer Activities, FY 96 Annual Report, Report No. SFIM-AEC-ET-CR-97013, pp. 155-156.

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#### Site Information:

- Aberdeen, MD
- Ogden, UT
- U.S Air Force Facility, Fort Worth, TX
- Milan Army Ammunition Plant (USAEC), TN
- DOE, Bear Creek, Oak Ridge National Laboratory, TN

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# Points of Contact:



# **Technology Specific Web Sites:**



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#### **Vendor Information:**

A list of vendors offering In Situ Biological Treatment is available from the Vendor Information System for Innovative Treatment Technologies (VISITT) developed by U.S. Environmental Protection Agency (EPA).



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# Health and Safety:

To be added

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# 4.5 Phytoremediation

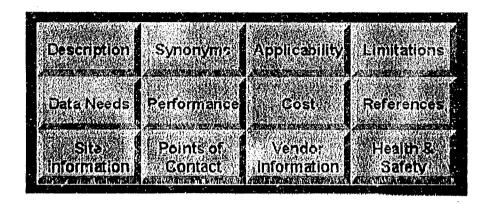
(In Situ Soil Remediation Technology)

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Technology	Description
Soil, Sediment, and Slu	ıdge
3.1 In Situ Biological T	<u>'reatment</u>
4.5 Phytoremediation	Phytoremediation is a process that uses plants to remove, transfer, stabilize, and destroy contaminants in soil and sediment. Contaminants may be either organic or inorganic.

#### **Description:**

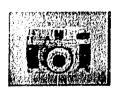


Figure 4-5: Typical In Situ Phytoremediation System

Phytoremediation is a process that uses plants to remove, transfer, stabilize, and destroy contaminants in soil and sediment. The mechanisms of phytoremediation include enhanced rhizosphere biodegradation, phyto-extraction (also called phyto-accumulation), phyto-degradation, and phytostabilization.

# ➤ Enhanced Rhizosphere Biodegradation

Enhanced rhizosphere biodegradation takes place in the soil immediately surrounding plant roots. Natural substances released by plant roots supply nutrients to microorganisms, which enhances their biological activities. Plant roots also loosen the soil and then die, leaving paths for transport of water and aeration. This process tends to pull water to the surface zone and dry the lower saturated zones.

# > Phyto-accumulation

Phyto-accumulation is the uptake of contaminants by plant roots and the translocation/accumulation (phytoextraction) of contaminants into plant

shoots and leaves.

## > Phyto-degradation

Phyto-degradation is the metabolism of contaminants within plant tissues. Plants produce enzymes, such as dehalogenase and oxygenase, that help catalyze degradation. Investigations are proceeding to determine if both aromatic and chlorinated aliphatic compounds are amenable to phyto-degradation.

## > Phyto-stabilization

Phyto-stabilization is the phenomenon of production of chemical compounds by plant to immobilize contaminants at the interface of roots and soil.

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Synonyms:

Vegetation-enhanced bioremediation.

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Applicability:

Phytoremediation may be applicable for the remediation of metals, pesticides, solvents, explosives, crude oil, PAHs, and landfill leachates.

Some plant species have the ability to store metals in their roots. They can be transplanted to sites to filter metals from wastewater. As the roots become saturated with metal contaminants, they can be harvested.

Hyper-accumulator plants may be able to remove and store significant amount of metallic contaminants.

Currently, trees are under investigation to determine their ability to remove organic contaminants from ground water, translocate and transpiration, and possibly metabolize them either to CO<sub>2</sub> or plant tissue.

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Limitations:

There are a number of limitations to phytoremediation in soil.

- The depth of the treatment zone is determined by plants used in phytoremediation. In most cases, it is limited to shallow soils.
- High concentrations of hazardous materials can be toxic to plants.
- It involves the same mass transfer limitations as other biotreatments.
- It may be seasonal, depending on location.
- It can transfer contamination across media, e.g., from soil to air.
- It is not effective for strongly sorbed (e.g., BCBs) and weakly sorbed contaminants.
- The toxicity and bioavailability of biodegradation products is not always known.
- Products may be mobilized into ground water or bioaccumulated in animals.
- It is still in the demonstration stage.

## • It is unfamiliar to regulators.

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Data Needs:

A detailed discussion of these data elements is provided in <u>Subsection 2.2.1</u> (Data Requirements for Soil, Sediment, and Sludge). In addition, detailed information is needed to determine the kinds of soil used for phytoremediation projects. Water movement, reductive oxygen concentrations, root growth, and root structure all affect the growth of plants and should be considered when implementing phytoremediation.

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Performance Data:

Currently, the Superfund Innovative Technology Evaluation (SITE) Program is attempting to demonstrate and evaluate the efficacy and cost of phytoremediation in the field at sites in Oregon, Utah, Texas, and Ohio.

USAEC is also leading the team of experts from EPA, Tennessee Valley Authority (TVA) and the Waterways Experimental Station (WES) to successfully demonstrate phytoremediation of explosive contaminated sites in Milan Army Ammunition Plant in Milan, TN.

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Cost:

US AEC estimated that the cost for phytoremediation of one acre of lead-contaminated soil to a depth of 50 cm was \$60,000 to \$100,000, whereas excavating and landfilling the same soil volume was \$400,000 to \$1,700,000.

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References:

Boyajian, G. E. and Devedjian, D. L., 1997. "Phytoremediation: It Grows on You", Soil & Groundwater Cleanup, February/March, pp. 22-26.

EPA, 1998. A Citizen's Guide to Phytoremediation, Technology Fact Sheet, EPA NCEPI, EPA/542/F-98/011.

EPA, 1996. A Citizen's Guide to Phytoremediation, Technology Fact Sheet, EPA NCEPI, EPA/542/F-96/014.

EPA, 1996. Recent Developments for In Situ Treatment of Metal Contamir sted Soils, EPA/542/R-96/008.

Schnoor, J.L., L.A. Licht, S.C. McCutcheon, N.L. Wolfe, and L.H. Carreira. 1995. "Phytoremediation of organic and nutrient contaminants," *Environ. Sci. Technol.* 29:318A-323A.

USAEC, 1997. "Phytoremediation of Lead" in Innovative Technology Demonstration, Evaluation and Transfer Activities, FY 96 Annual Report, Report No. SFIM-AEC-ET-CR-97013, pp. 89-92.

U.S. DOE, 1995. "Bioremediation of High Explosives by Plants," in *Technology Catalogue, Second Edition*, Office of Environmental Management Office of Technology Development, DOE/EM-0235, pp. 169-

172.

A comprehensive list of 850 references on phytoremediation are available at Remediation Technologies Development Forum (RTDF) Phytoremediation Action Team Web Site. Click to access



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#### Site Information:

- McCormick & Baxter SUPERFUND Site, Portland, OR (Wood treatment site)
- Argonne National Laboratory
- Craney Island Fuel Terminal, U.S. Navy, Portsmouth, VA
- EPA S.I.T.E. Program, Ogden, Utah
- Ohio (Former metal plating site)
- DOE Demo: Savannah River Site, SC
- DOE Savannah River Site, SC

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#### **Points of Contact:**



#### **Technology Specific Web Sites:**



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#### **Vendor Information:**

A list of vendors offering In Situ Biological Soil Treatment is available from the Vendor Information System for Innovative Treatment Technologies (VISITT) developed by U.S. Environmental Protection Agency (EPA).



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# APPENDIX D

Preliminary Cost Estimate CKD landfill Cap





Preston Engineering Inc. 4436 N. Brady Street Davenport, Iowa

ATTN: Mr. Morris Preston, P.E.

RE: GSI No. 993181

**Preliminary Construction Cost Estimate** 

CKD Landfill Cap

**Dixon Marquette Cement Company** 

Dixon, Illinois 61021

Dear Mr. Preston:

Geotechnical Services, Inc. (GSI) is pleased to provide this preliminary cost estimate for the construction of the CKD Landfill Cap. As requested in our telephone conversation of June 25, 1999, and June 28, 1999 the following information presents our preliminary cost estimate for the construction of a three (3) foot thick low permeability cap with hydroseeded vegetative layer for the Cement Kiln Dust (CKD) Landfill to be located at the Dixon Marquette Cement Company in Dixon, Illinois.

For estimating purposes the following data has been utilized:

- 1. Landfill cap: 24 acre area, capped with a 3 foot recompacted low permeability layer without drainage layer. A six inch vegetative layer will be placed over the recompacted low permeability cap. Hydroseeding will be utilized to seed the vegetative layer.
- 2. On site sandy clay soil will be amended and utilized for the cap borrow material, borrow transport distance will be 2,000 feet or less.
- 3. On site sandy clay borrow does not meet required permeability (k) of  $1 \times 10^{-7}$  cm/sec or less.
- 4. Prior to construction, geotechnical laboratory testing will be performed to identify optimum soil amendment quantities to achieve required permeability value based on Standard Proctor compaction criteria.
- 5. Construction QA/QC will be performed by an independent testing laboratory.
- 6. Bentonite soil amendment will be utilized to achieve recompacted soil layer permeability (k) of 1 x 10<sup>-7</sup> cm/sec or less for the cap.

GEOTECHNICAL, MATERIALS ENGINEERING & ENVIRONMENTAL CONSULTANTS

258 EAST 90TH STREET DAVENPORT, IOWA 52808-7341 (319) 285-8541 • FAX (319) 285-8545



- 7. On site water will be used for moisture conditioning at no charge to the contractor.
- 8. On site sandy clay borrow will be mixed/amended at a rate of approximately 2 pounds bentonite /cu.ft. of borrow, moisture conditioned in 8 inch loose lifts and compacted with sheeps foot roller with appropriate lift bending.
- 9. Construction equipment consisting of CMI or Caterpillar recycler/pulverizer, tank truck with spray bar, earth moving pans, grader, and sheeps foot compactor self propelled.
- 10. Off site contractor will provide equipment operators and manual labor for the addition of the bentonite amendment and construction of the cap.
- 11. Bentonite amendment to be supplied by Bentonite Performance Minerals, delivered by truck, 100# bags on pallets.
- 12. Estimated construction completion: 4 months.

Based on the above information, the following preliminary cost estimate for the landfill cap has been developed:

Item	Material (Per Acre)	Labor (Per Acre)	Cost/Acre (M & L)	Total Cost
Hydroseed	\$ 900	\$ 600	\$ 1,500	\$ 36,000
Vegetative Layer, 6"	\$ 9,700	\$ 2,400	\$12,100	\$290,400
Bentonite Amendment	\$15,700	\$ 2,100	\$17,800	\$427,200
Haul/Place Borrow		\$ 8,700	\$ 8,700	\$208,800
Mixing & Compaction		\$21,000	\$21,000	\$504,000
Construction QA/QC		\$ 3,000	\$ 3,000	\$ 72,000
Amendment Lab Testing		\$ 200	\$ 200	\$ 4,800



Total Estimated Cost \$1,543,200

Estimated Cost/Acre \$ 64,300

Total Estimated Cost (w/20% Contingency) \$1,851,840

If you have any questions or require additional information concerning the above information, material and labor rates and the estimated cost for the construction of this landfill cap, please call me at 319-285-8541.

Sincerely,

GEOTECHNICAL SERVICES, INC.

George D. Komisarek, P.E., P.G.

Vice President

cc: file

Mr. Neil DeRynck Dixon Marquette Cement Company 1914 White Oak Lane Dixon, Illinois 61021

Kernivaru

# APPENDIX E

Leachate Collection System

Dixon Marquette Cement 98-516

Proposed Leachate Collection System January 27, 7000

The CKD landfill at Disca Marquette Cement generates leachate that enters the bedrock beneath the landfill and travels downgradient toward the Rock River. To meet State of Illinois requirements, a leachate collection system must be installed.

The CKD directly overlies the Pecatonica Formation, a highly jointed dolomite about 35 feet thick. The joints in the rock are essentially vertical, and run in two preferred directions, roughly north and south, and east to west. Groundwater flow beneath the landfill is probably primarily along the joints. Consequently, any effective and efficient leachate collection system should be designed to intercept groundwater in the joints.

Traditional dewatering systems use multiple vertical wells. This type of system would work as a leachate collection system at the site, but only if wells could be placed so that the screens would intersect the joints. Because the number of and locations of the joints are not known, well placement becomes a problem. One way to solve the problem is to install a large number of vertical wells in the hope that some percentage of them will coincide with joints. This becomes a trial-and-error method that potentially wastes a lot of money in exploratory drilling.

The geologic conditions are very suitable, on the other hand, for horizontal wells. A single horizontal well, drilled at some angle to the planes of the joints, will intersect the joints. Each horizontal well can replace dozens of vertical wells. There are cost savings that come with using horizontal wells. Vertical wells would need a pump placed in each well, with electricity delivered to each pump. Surface piping connecting the wells would need to be trenched in for hundreds of feet. The multiplicity of recovery pumps, and the cost of maintaining them, would be many times more than the single pump in a horizontal well. Electrical costs in operating a single well would also be less. The saving in operation and maintenance increases as the duration of the project increases.

Unless a large number (hundreds?) of vertical wells are drilled, the probability of failure of vertical wells to capture the leachate is high. The recommended collection system, therefore, is one that uses a horizontal well.

The proposed system consists of one horizontal well. It will run east of and parallel to the haul road on the west side of the landfill. It will be about 8 to 10 inches in diameter, and be located near the base of the Pecatonica Formation. This location should allow the well to capture a large majority of the groundwater in the uppermost aquifer that has traveled beneath the landfill. It should also dewater the Pecatonica enough so that groundwater is no longer in contact with CKD at the base of the landfill.

The horizontal well could be drilled from either the north or south end. The drill stern will enter the ground at a shallow angle and be rotated so it becomes horizontal at a depth of about 30 feet. At the other end, the drill stem will again be rotated, this time upward, to exit at a point about 2000 feet from the entry. A well screen will then be pulled into the hole as the drill is backed out.

A single large capacity submersible pump will be used in the well to recover groundwater, which will be transferred to a holding basin before treatment and discharge to the Rock River.

# PRESTON ENGINEERING, INC.

PROJECT ESTIMATE for:	Leachate system, 1 horizontal well
Project #:	98-516
Location:	Dixon Marquette
Date of Estimate:	19-Dec 00

DESCRIPTION	TIAU YTITHAUQ	RATE	TOTAL	
PERSONNEL				
Principal	18 hour	\$100.00	\$1,600	management
Principal	16 hour	\$100.00	\$1,600	4 site visits
Level 3 Scientist	40 hour	\$70.00	\$2,800	design
Level 3 Scientist	130 hour	\$76.00	\$9,100	travel, 54 trips
Level 3 Scientist	192 hour	\$70.00	\$13,440	install well
Level 3 Scientist	0 hour	\$70.00	\$0	manifolding, trenching
Level 3 Scientist	40 เ.วเศ	\$70.00	\$2,800	startup
Level 3 Scientist	0 hour	\$70.00	\$0	
Level 3 Scientist	0 hour	\$70.00	\$0	
Level 3 Scientist	0 hour	\$70.00	\$0	
Level 2 Technician	8 hour	\$49.00	\$392	
TDAY (5)				\$31,732
TRAVEL Mileage	7560 mile	*0.00	*0.070	
Tolis, parking	7560 mile 54 each	\$0.38 \$0.60	\$2,873	
Per Diem	0 day	\$40.00	\$32 \$0	
Lodging	•	*	•-	
cooging	0 day	\$50.00	\$0	
EQUIPMENT AND SUPPLIES				
Surveying equipment	1 day	\$25.60	\$25	
Submersible pumps and contn	1 each	\$10,000.00	\$10,000	
Equipment sheds	1 each	\$3,000.00	\$3,000	
Computer drafting	4 hour	\$10.00	\$40	
DRILLING SUBCONTRACTORS				\$15,970
Mobilization	1 each	\$500.00	\$500	
Bedrock drilling	2000 feet	\$200.00	\$400,000	
Well installation (PVC)	2000 feet	\$50.00	\$160,000	
Flush mount covers	0 each	\$125.0G	\$100,000	
Well development	10 hour	\$125.00 \$125.00	\$1,250	
Decontamination	0 hour	\$125.00		
Per diern	23 day	\$150.00	\$0 \$3,450	
	,	• • • • • • • • • • • • • • • • • • • •	0-7.00	
EXCAVATOR SUBCONTRACTOR				
Backhoe	0 hour	\$75.00	\$0	
ELECTRICAL SUBCONTRACTOR				
Connect pump	16 hour	\$75.00	\$1,200	
SUBCONTRACTOR HANDLING	10 percent	\$506,400	\$50,640	
CONTINGENCY	15 percent	\$604,749	\$90.711	
TOTAL		<u></u>	\$595,454	1

## PRESTON ENGINEERING, INC.

PROJECT ESTIMATE for:

Project #:

Location:

Date of Estimate:

Leachate system, 100 vertical wells

Project #:

Dixon Marquette

28-Jan-00

				•
DESCRIPTION	QUANTITY UNIT	RATE	IQTAL	
PERSONNEL				
Principal	16 hour	\$100.00	-	management
Principal	16 hour	\$100.00	\$1,600	4 site visits
Level 3 Scientist	40 hour	\$70.00	\$2,800	design
Level 3 Scientist	130 hour	\$70.00	\$9,100	travel, 54 trips
Level 3 Scientist	272 hour	\$70.00		install wells
Level 3 Scientist	160 hour	\$70.00	\$11,200	manifolding, trenching
Level 3 Scientist	80 hour	\$70.00	. \$5,600	startup
Leval 3 Scientist	O hour	\$70.00	\$0	
Level 3 Scientist	O hour	\$70.00	\$0	
Level 3 Scientist	O hour	\$70.00	\$0	
Level 2 Technician	8 hour	\$49.00	\$392	
				\$51,332
TRAVEL Mileage	7560 mile	\$0.38	\$2,873	
-	54 each	\$0.60	\$32	•
Tolis, parking	O day	\$40.00	\$0	
Per Diem	O day	\$50.00	\$O	
Lodging .	O day	1110.00	***	
EQUIPMENT AND SUPPLIES				
Surveying equipment	1 day	\$25.00	\$25	
Submersible pumps and controll	100 each	\$3,000.00	\$300,000	
Equipment sheds	3 each	\$3,000.00	\$9,000	
Computer drafting	8 hour	\$10.00	\$80	\$312,010
DRILLING SUBCONTRACTORS				4512,010
Mobilization	1 each	\$500.00	<b>\$500</b>	GSI
Bedrock drilling	3000 feet	\$25.00	\$75,000	
Well installation (PVC)	3000 feet	\$11.00	\$33,000	
Flush mount covers	100 each	\$125.00	\$12,500	
Well development	50 hour	\$125.00	\$6,250	
Decontamination	O hour	\$100.00	\$0	
Per diem	34 day	\$150.00	\$5,100	
EXCAVATOR SUBCONTRACTOR				
Backhos	80 hour	\$75.00	\$6,000	
ELECTRICAL SUBCONTRACTOR				
Connect 100 pumps	200 hour	\$75.00	\$15,000	
SUBCONTRACTOR HANDLING	10 percent	\$153,350	\$15,335	
CONTINGENCY	15 percent	\$532,027	\$79.804	
TOTAL			<b>#811,831</b>	]

# REQUEST FOR ADJUSTED GROUNDWATER STANDARDS &

DIXON MARQUETTE CEMENT COMPANY DIXON, ILLINOIS

PREPARED BY PRESTON ENGINEERING, INC 4436 NORTH BRADY AVENUE DAVENPORT, IOWA 52806 TELEPHONE: (319) 388-8288 FAX: (319) 388-9003

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# **ATTACHMENTS**

Summary Tables
Background Tables
Downgradient Tables
Graphs

#### I. INTRODUCTION

Groundwater samples have been collected from wells around the Dixon Marquette Cement (DMC) landfill for several years. This information from wells (MW5-S, MW6-S, MW7-S, and MW8-S) at the boundary of the proposed zone of attenuation (ZOA) indicate values for some chemical parameters that exceed the background concentration established for the background wells (MW 1-WT, MW 1-S). A statistical analysis was performed on the data covering thirty-nine (39) parameters.

The background concentration was evaluated by establishing the 99% confidence level for each parameter. Levels in downgradient wells that are higher than this confidence interval indicates a statistically significant increase over background.

The concentration of parameters in downgradient wells was examined using statistical methods. The 99.5% confidence level was established for each parameter. A trend analysis was also performed to determine if the level in each well appeared to be increasing or decreasing over time. The data, trend line, and 99.5% confidence level were plotted for each well and each parameter.

A narrative discussion of each parameter was provided.

Groundwater criteria were established for each parameter. Levels were selected that could be met at monitoring wells MW 5-S, MW 6-S, MW 7-S and MW 8-S. The criteria selected was the higher of the Section 620 groundwater standards or the 99.5% confidence interval. Higher levels were set for some parameters where it appeared there was an increasing trend and where levels where higher at MW 3-S which is upgradient. In these cases the 99.5% confidence level for MW 3-S was used as a criteria.

## II. BACKGROUND CONCENTRATIONS

Statistical analysis was performed on three background monitoring wells, MW 1-WT, 1-S, and 1-D. Six methods were used to calculate the upper limit concentration to compare the downgradient monitoring wells with the background.

- Student's T Distribution: When the parameter was detected in all of the sample results, this method was used to calculate the upper limit concentration. The student's t distribution is used for a small data set, less than 30 samples, and assumes that the sample set is normal.
- Detection Limit Divided by 2 (DL/2): In cases where 15 percent or less of the data was below the
  detection limit, the values below the detection limit were replaced with one-half of the detection
  limit. Student's t test was then applied.
- Cohen's Method: When less than 50% of the samples were reported as below the detection limit, the upper background limit concentration was calculated using this method. Cohen's method provides estimates of the mean and standard deviation when there are observations less than the detection limit. The method assumes that the sample set is normally distributed and the detection limit remains the same through the sampling period.
- . Test of Proportions: When more than 50% of the samples were reported as below the detection limit, a test of proportions was performed on the downgradient monitoring well. This procedure does not establish a background concentration since it is a comparison of two sampling sets. Using the sample sizes and the number of samples that were below the detection limit in each sample set, the test indicates if the downgradient concentration is statistically different than the background concentrations.
- Lowest Non Detect: When the concentration was below detection limits during the entire sampling period, the lowest non-detect was used as the upper limit.
- Method Detection Limit: The method detection limit was established as the upper limit for
  volatile and semi-volatile compounds. However, any detection is considered to be an exceedance
  of the background concentration.

Using data from only one monitoring well, the 99<sup>th</sup> percentile was calculated for the upper limit in the student's t distribution and the Cohen methods. The two sided 99<sup>th</sup> percentile confidence interval was used to calculate background levels for pH. This was performed in accordance to Title 35 of Illinois Administrative Code, Section 811.320.e, entitled: "Statistical Analysis of Groundwater Monitoring Data."

The sampling results of MW 1-S, 1-WT, and 1-D, and the statistical analysis is summarized in the section entitled "BACKGROUND TABLES."

#### III. STATISTICAL APPROACH OF DOWNGRADIENT MONITORING WELLS

The arithmetic mean and standard deviation were calculated for each parameter and each well. When the result was reported as below the detection limit, the detection limit was used in the calculations. These results are denoted in the tables by italicizing the value.

The data set was compared to the average plus one standard deviation and the average plus two standard deviations. Those that were greater than the average plus one standard deviation were represented with bold characters in the tables. The results that were greater than the average plus two standard deviations were highlighted in the tables.

The one sided 99.5% confidence interval was calculated for each parameter and each of the four downgradient monitoring wells (5-S through 8-S). For pH, the two-sided 99% confidence interval was calculated. Since there were less than 30 sampling events, the student's t distribution was used to calculate the interval. The four intervals were compared, and the maximum value was chosen to represent the results of the four downgradient monitoring wells.

Sampling data and statistical analysis for each well is located in the section of this document entitled "DOWNGRADIENT TABLES." Summaries of the confidence intervals, background concentrations, and the Class I Groundwater Standards are found in the section entitled "SUMMARY TABLES."

The data was graphed with the 99.5% confidence interval and the Class I Groundwater Standard, when applicable. The standards are published in Title 35 of the Illinois Administrative Code Section 620.410 entitled Groundwater Quality Standards for Class I: Potable Resource Groundwater. A trend line was added to each of the wells for each parameter. Since the results did not indicate a clear trend, a linear trend or the best-fit line was chosen to represent the data.

The graphs of each parameter are found in the section entitled "GRAPHS" of this document,

## IV. GROUNDWATER IMPACT- INORGANIC COMPOUNDS

In the following sections, each inorganic compound in MW 5-S through 8-S is discussed with respect to the background concentrations. When comparing these wells to background concentrations, both MW 1-WT and 1-S were used since the groundwater is shallow near White Oak Lane. Since MW 1-D monitors the upgradient groundwater in the Saint Peter Formation, the levels were not compared to the MW 5-S through 8-S.

#### Aluminum

The upper limit in the background MW 1-WT was calculated to be 0.454 mg/L using Cohen's method. The upper limit in MW 1-S was set at 0.04 mg/L, the lowest non-detectable limit. All but three results were greater than the MW 1-S background limit. All concentrations in MW 5-S were below the background concentration in 1-WT. In MW 6-S, two values, 1.74 and 3.01 mg/L, were significantly higher than the 1-WT background concentration. There were two results in MW 7-S that were slightly greater than the 1-WT background. These were 0.88 mg/L on July 26, 1999 and 0.662 on March 2, 2000. In the July and November 1999 samplings, the results in the MW 8-S were also slightly greater than the 1-WT background. These values were 0.51 and 0.665 mg/L, respectively.

The average concentrations of aluminum in the four downgradient monitoring wells were below the background concentration in MW 1-WT. Each, however, were greater than upper limit of 0.04 mg/L in MW 1-S.

## Antimony

Since antimony was not detected in the background well 1-WT, the upper limit was set to be 0.001 mg/L, the lowest not detectable limit. Antimony was detected only once during the monitoring period in one of the wells. The parameter was detected in MW 7-S at 0.003 mg/L on March 2, 2000.

To determine if there is a difference between the background MW 1-S and the four downgradient wells, a test of proportions was performed. The test of proportions indicates that the result was not different than the background MW 1-S. Therefore, the landfill has had very little affect on this parameter in the groundwater.

#### Arsenic

The upper limit background concentration was calculated to be 0.003 in MW 1-WT using the student's t distribution. The concentration of arsenic ranged from 0.001 to 0.005 mg/L in MW 5-S, 0.001 to 0.004 in MW 6-S, 0.0017 to 0.044 mg/L in MW 7-S and 0.001 to 0.005 mg/L in MW 8-S. A total of four sample results, one from each well, were greater than the upper limit concentration in MW 1-WT.

A test of proportions was performed to indicate whether or not the arsenic concentrations in the downgradient wells were significantly different than the concentrations in MW 1-S. Although arsenic was detected in all four of the downgradient monitoring wells, the concentrations were not considered to be significantly different than MW 1-S except for MW 5-S. The landfill has had very little affect on this parameter in the groundwater.

#### Barium

The background concentration of barium was calculated to be 0.227 and 0.091 mg/L in the MW 1-WT and 1-S, respectively. These limits were calculated using the student's t 99% confidence interval. Barium was detected in all four of the downgradient monitoring wells. The concentrations ranged from 0.025 to 0.35 mg/L. Four of the samples were greater than 0.091 mg/L in MW 5-S, however only 1 was greater than 0.227 mg/L. All seven results were greater than the MW 1-S background limit in MW 6-S. Two of the results were greater than the 1-WT background limit. There were no results that were greater the background concentrations in MW 7-S, and only 1 that was greater than the MW 1-WT background concentration in MW 8-S.

The average concentration of barium in MW 5-S through 8-S was 0.138, 0.258, 0.042, and 0.075 mg/L, respectively. Monitoring well 6-S was the only average that was greater than the background concentration of 0.227 mg/L. Therefore, the landfill has had very little affect on this parameter in the groundwater.

#### Beryllium

Since beryllium was not detected in the background monitoring wells, 1-S or 1-WT, the upper background limit is 0.001 mg/L, the lowest non-detectable limit. On March 2, 2000, beryllium was detected in MW 8-S. This was the only time that the parameter was detected in the four shallow wells along the Rock River. Therefore, the landfill has had very little affect on this parameter in the groundwater.

#### Boron

To determine if the downgradient wells are significantly different than the background monitoring wells, a test of proportions was performed. The tests indicated that only MW 8-S was significantly different than the background MW 1-WT. All of the wells were significantly different than MW 1-S.

The concentration of boron ranged from 0.05 to 1.95 mg/L in the four downgradient monitoring wells. The average concentration of boron in MW 5-S through 8-S was 0.464, 0.424, 0.436, and 0.208 mg/L, respectively. This parameter in the groundwater has been significantly impacted by the landfill.

## Cadmium

Since cadmium was not detected in the background monitoring wells 1-WT or 1-S, the upper background limit is 0.001 mg/L, the lowest non-detectable limit for each. Cadmium was detected in MW 5-S, 6-S, and 8-S during the March 2000 sampling event only. The average concentration for each of the four downgradient monitoring wells was 0.001 mg/L, the detection limit. Therefore, the landfill has had very little affect on the concentrations of cadmium in the groundwater.

## Calcium

The background concentration of calcium was calculated to be 109 mg/L in the MW 1-S and 126 mg/L in MW 1-WT using the upper limit of the student's t 99% confidence interval. The concentration of calcium in monitoring wells 5-S, 6-S, and 8-S were consistently greater than the background limits during the entire sampling period. The concentrations ranged from 24.6 to 345 mg/L. The average concentrations of calcium in these three wells were 240, 234, and 237 mg/L, respectively. Calcium concentrations in MW 7-S were consistently below the background concentrations ranging from 11.9 to 43.4 mg/L with an average of 35 mg/L. The concentration of calcium in the groundwater has been significantly impacted by the landfill.

#### Chromium

The background concentrations were calculated to be 0.002 mg/L in both MW 1-WT and 1-S using Cohen's method. Chromium was detected in all four of the downgradient monitoring wells 5-S through 8-S. The maximum concentration was found in MW 8-S during the March 2000 sampling event. The concentration was found to be 0.04 mg/L. The average concentration of chromium in MW 5-S through 7-S was found to be 0.003 mg/L. The average concentration in MW 8-S was calculated to be 0.010 mg/L. Concentrations of chromium in the groundwater have been slightly affected by the landfill.

### Cobalt

Using Cohen's method, the upper limit of the 99% confidence interval was calculated to be 0.003 mg/L in MW 1-S. In MW 1-WT, the background limit was calculated to be 0.008 mg/L using the student's t distribution. Cobalt was detected in all four of the downgradient monitoring wells 5-S through 8-S during the sampling period. Most of the concentrations were below the 0.008 mg/L background level. Three of the exceedances were in MW 6-S. In the March 2000 sampling event, MW 8-S also exceed the background concentration, having a result of 0.035 mg/L.

The average concentration of cobalt in MW 5-S through 8-S is 0.002, 0.012, 0.001, and 0.006 mg/L. The landfill has had a small affect on the concentration of cobalt in the groundwater.

Copper

The background concentration was calculated to be 0.012 mg/L in MW 1-WT using student's t distribution and 0.008 mg/L with Cohen's method in MW 1-S. Copper was detected in all four of the monitoring wells. The concentrations ranged from below the detection limit to 0.113 mg/L. In MW 5-S there was one result that was greater than the 1-WT background limit. The rest were below 0.008 mg/L. In MW 6-S, there were three results that were greater than 0.012 mg/L and three that were below 0.008 mg/L. All of the sampling results were less than or equal to 0.008 mg/L in MW 7-S. In MW 8-S, there were three results greater than 0.012 mg/L and one that was greater than 0.008 mg/L.

The average concentration of copper in the four downgradient monitoring wells 5-S through 8-S is 0.006, 0.013, 0.005, and 0.024 mg/L, respectively. The concentration of copper has been affected slightly due to the landfill.

#### Lead

Since lead was not detected in the background MW 1-S, the upper background limit is 0.001 mg/L, the lowest non-detectable limit. In MW 1-WT, the background concentration was calculated to be 0.006 mg/L. In MW 5-S, lead was detected in two of the samples and was reported at the detection limit. The average lead concentration in this monitoring well was calculated to be 0.001 mg/L. In MW 6-S, there were two sample events that had detectable concentrations. The concentration of these detected concentrations were 0.015 and 0.008 mg/L and were greater than the Class I Groundwater Standard of 0.0075 mg/L. The average concentration was calculated to be 0.004 mg/L. Four out of the seven sample sets in MW 7-S had levels that were detected. These were less than 0.006 mg/L. In MW 8-S, five of the sample had levels of lead that were detectable. The concentrations were below 0.006 mg/L.

Of all of the sampling events, two sample results in MW 6-S were greater than the MW 1-S upper limit background concentration. These concentrations were also greater than the Class I Groundwater Standard. Lead concentrations in the downgradient monitoring wells has been slightly affected by the landfill.

#### Iron

The background concentration was calculated to be 0.465 mg/L in MW 1-S using Cohen's method and 0.891 in MW 1-WT using student's t distribution. Iron was detected in all of the sample locations during all of the sample events. The average concentrations were calculated to be 0.898, 2.665, 0.805, and 1.654 for MW 5-S through 8-S, respectively. The concentration of iron in the groundwater has been significantly impacted by the landfill.

Magnesium

The background concentration was calculated to be 53.9 and 50.5 mg/L in the MW 1-S and 1-WT, respectively. Both of these confidence intervals were calculated using the upper limit of the student's t distribution. The average concentrations were calculated to be 82.3, 55.16, 9.3, and 100.1 mg/L for monitoring wells 5-S through 8-S, respectively. Approximately 72% of the samples were greater than the MW 1-WT background concentration in MW 5-S. Four of the seven sampling events had concentrations greater than 50.478 mg/L. The concentration of magnesium was greater than 50.478 mg/L in MW 8-S. In MW 7-S, the greatest concentration was 15.1 mg/L. The concentration of magnesium in the groundwater has been significantly impacted by the landfill.

Manganese

The background concentration was calculated to be 0.045 mg/L in the MW 1-S and 2.687 mg/L in MW 1-WT using the upper limit of the student's t 99% confidence interval. Two concentrations in the four downgradient wells had concentrations that were greater than 2.687 mg/L. Monitoring well 6-S had a manganese concentration of 3.11 mg/L during the July 1999 sampling event. Monitoring well 8-S had a manganese concentration of 4.3 mg/L during the March 2000 sampling event. The average concentrations were 0.229, 1.695, 0.094, and 0.763 in the four respective monitoring wells, 5-S through 8-S. Therefore, the landfill has had very little affect on the concentration of manganese in the groundwater.

Mercury

Since mercury was not detected in the background MW 1-S, the upper background limit is 0.0002 mg/L, the lowest non-detectable limit. A test of proportions showed that the wells were not significantly Mercury was not detected in the four shallow wells near the Rock River during the monitoring period. There has been very little affect on the concentration of mercury from the landfill.

Nickel

The background concentration was calculated to be 0.007 mg/L in MW 1-S using Cohen's method. In MW 1-WT the 99% confidence interval was calculated to be 0.023 mg/L using student's t distribution. There were three sampling results that were greater than the 1-WT background concentration in the four downgradient monitoring wells 5-S through 8-S. Two concentrations, 0.033 and 0.038 mg/L, were in the sampling data from MW 6-S. During the March 2000 sampling event, MW 8-S had a nickel concentration of 0.058 mg/L. The average concentration of nickel in MW 5-S through 8-S were 0.005, 0.020, 0.010, and 0.014 mg/L, respectively. The landfill has had a small affect on the concentration of nickel in the groundwater.

#### Potassium

Since the data had 50% of non-detects in the background well, MW 1-S, a test of proportions was performed to determine if the downgradient wells were significantly different. Each of the four downgradient wells had levels of potassium that was detected during all of the seven sampling events. The test of proportions proved that the downgradient and upgradient wells were significantly different.

The upper limit of the 99 percentile background concentration in MW 1-WT was calculated to be 4.0 mg/L, using the student's t distribution. Each of the four downgradient wells was significantly higher than the background. The average concentrations in monitoring wells 5-S and 6-S were calculated to be 86.0, and 55.0 mg/L. Monitoring wells 7-S and 8-S were higher with averages of 632.6 and 245 mg/L, respectively. The concentration of potassium in the groundwater has been significantly impacted by the landfill.

## Selenium

The 99<sup>th</sup> percentile was calculated to be 0.004 mg/L in background well 1-S. In MW 5-S, concentrations ranged from below detection limits to 0.006 mg/L. The average concentration was 0.002 mg/L in this well. Monitoring well 6-S and MW 7-S had an average selenium concentration of 0.003 and 0.002 mg/L, respectively. The range for each of these wells was from below detection limits to 0.005 mg/L. Monitoring well 8-S was significantly different than the background monitoring well. The average concentration in this well was calculated to be 0.005 mg/L and the data ranged from 0.001 to 0.014 mg/L.

A test of proportions was performed to determine if the downgradient wells were considered to be significantly different than the background well 1-WT. In the downgradient monitoring wells 5-S through 8-S, there were 6 to 7 samples that had levels of selenium that was detected. The test of proportions confirmed that the downgradient wells were significantly different than MW 1-WT. Comparing both background concentrations to the four downgradient monitoring wells, the landfill has affected the concentration of selenium.

#### Sodium

The background concentration was calculated to be 12.8 mg/L in the MW 1-S using the upper limit of the student's t 99% confidence interval. In MW 1-WT, the 99th percentile was calculated to be 53.7 mg/L. In MW 5-S, the concentrations ranged from 9 to 78.5 mg/L with an average sodium concentration of 37.9 mg/L. The average concentration of sodium in MW 6-S was calculated to be 46.4 mg/L. The sample results ranged from 27.3 to 64.9 mg/L. In MW 7-S, the sampling results ranged from 25.8 to 49.9 mg/L with an average of 40.9 mg/L. In MW 8-S, the average was calculated to be 39.7 mg/L and the sample result ranged from 8.2 to 52.4 mg/L. Sodium concentration in the groundwater has been significantly affected by the landfill.

#### Thallium

To determine if the downgradient monitoring results were significantly different than MW 1-S and MW 1-WT, tests of proportions were performed. The results of these tests indicated that there was no significant difference between the background, MW 1-S and MW 1-WT, and the downgradient monitoring wells, 5-S through 8-S. Thallium was detected once during the monitoring period in MW 6-S at the detection limit, 0.001 mg/L. In MW 8-S, there were two sampling events in which the parameter was detected. The results were 0.002 and 0.005 mg/L. The average thallium concentration for this monitoring well was calculated to be 0.002 mg/L. The landfill has had very little affect on this parameter in the groundwater.

#### Vanadium

The upper limit of the 99<sup>th</sup> percentile was found to be 0.007 mg/L in the background MW 1-WT using the student's t distribution. In MW 5-S, there was one concentration that exceeded the background. The concentration was reported during the March 2000 sampling event and had a concentration of 0.023 mg/L. The average vanadium concentration was calculated to be 0.007 mg/L. Results of MW 6-S were similar to MW 5-S. There was only one sample result, 0.023 mg/L, that was greater than the background concentration. The average vanadium concentration for this well is 0.008 mg/L. The sampling results in MW 7-S ranged from 0.006 to 0.022 mg/L. The average concentration was calculated to be 0.013 mg/L. All of the results were less than the background concentration in MW 8-S. The average concentration was calculated to be 0.003 mg/L.

A test of proportions was performed on the data to determine if the results were significantly different than the background MW 1-S. The test of proportions found that only MW 7-S is considered to be different than the results of the background MW 1-S. The landfill has had a small impact affect on the concentration of vanadium in the groundwater.

## Zinc

The background concentration was calculated to be 0.058 and 0.014 mg/L in the MW 1-S and 1-WT, respectively. Each of these concentrations was calculated using the upper limit of the student's t 99% confidence interval. There was one sample result in the four monitoring wells, 5-S through 8-S that was greater than the background concentration of 0.058 mg/L. The concentration, 0.363 mg/L. was reported in the March 2000 sampling event in MW 8-S. The average zinc concentrations for MW 5-S through 8-S are: 0.017, 0.015, 0.016, and 0.075, respectively. The landfill has had very little affect on this parameter in the groundwater.

## Total Alkalinity

The background concentration was calculated to be 253 and 342 mg/L in the MW 1-S and 1-WT, respectively Each of these concentrations was calculated using the upper limit of the student's t 99% confidence interval. In MW 5-S through 8-S, the alkalinity was consistently higher than the background concentrations. Only one value was less than the MW 1-WT background concentration at 324 mg/L in MW 5-S. The average concentrations for the four downgradient wells were 442, 389, 396, and 287 mg/L. The concentration of total alkalinity in the groundwater has been significantly impacted by the landfill.

## Phenol Alkalinity

Phenol alkalinity was not detected in the three samples of MW 1-WT, thus the upper limit of the background concentration was set at 1.0 mg/L, the lowest detection limit.

To determine if the downgradient monitoring wells were significantly different than the background concentrations in MW 1-S, a test of proportions was performed on each downgradient well. Monitoring well 7-S has the only detectable levels of phenol alkalinity. The test of proportions indicates that this is not significantly different than the background. However, the concentrations range from below the detection limit to 160 mg/L. The average concentration of phenol alkalinity was calculated to be 73.5 mg/L.

#### Bicarbonate Alkalinity

The background concentrations were calculated to be 250 and 334 mg/L in the MW 1-S and 1-WT using the upper limit of the student's t 99% confidence interval. Sample results ranged from 314 to 540 mg/L in MW 5-S. The average concentration was calculated to be 410 mg/L. In MW 6-S, results ranged in concentration from 325 to 410 mg/L. The average concentration in this well was calculated to be 366 mg/L. Bicarbonate alkalinity in MW 7-S ranged from 154 to 475 mg/L and had an average concentration of 366 mg/L. In MW 8-S, all of the results were less than the background concentration of 334,029 mg/L. The average was calculated to be 276 mg/L in MW 8-S. The concentration of bicarbonate alkalinity has been slightly impacted due to the landfill.

## Ammonia Nitrogen

Since there were more than 50% of samples that had non-detectable results in the background monitoring wells 1-S and 1-WT; a test of proportions was performed on each downgradient monitoring well. The test showed that there was no significant difference in concentrations between the both of the background concentrations and the concentrations in the four downgradient monitoring wells. There were a total of two sample results that had detectable levels of ammonia nitrogen in the downgradient wells. These were in MW 5-S and 7-S with values of 0.19 and 0.94 mg/L, respectively. Therefore, the landfill has had very little affect on this parameter in the groundwater.

#### Chloride

The background concentration was calculated to be 30.8 and 25.4 mg/L in MW 1-S and 1-WT, respectively, using the upper limit of the student's t 99% confidence interval. In MW 5-S, the concentration of chloride ranged from 21.6 to 125 mg/L with an average of 62.7 mg/L. In MW 6-S, the sampling results ranged from 71.1 to 246 mg/L. The average concentration of chloride was calculated to be 152.9 mg/L. The average chloride concentration in MW 7-S was found to be 82.6 mg/L, the data ranged from 48.8 to 110 mg/L. The concentration of chloride ranged from 113 to 198 in MW 8-S. The average was calculated to be 153 mg/L. The concentration of chloride in the groundwater has been significantly impacted by the landfill.

## Chemical Oxygen Demand (COD)

Since the chemical oxygen demand was not detected in the background MW 1-S, the upper background limit is 5.0 mg/L, the lowest non-detectable limit. When detected, the sample results were greater than 5 mg/L in each of the four downgradient monitoring wells.

A test of proportions was performed to determine if there was a significant difference between the COL concentrations in the background monitoring well, 1-WT. The background well had three of the seven samples that were detected. The concentrations were 44, 25.6, and 42.9 mg/L. Although, the test of proportions indicated that there was no difference in the data, it appears that there is a difference in three of the downgradient monitoring wells. The four detected concentrations in MW 5-S were 26, 57.9, 25.7, and 203 mg/L. Monitoring well 6-S had three detected concentrations. The results are 100, 123, and 94.1 mg/L. Two sample events had quantitative results in MW 8-S. The results were 30.3 and 158 mg/L. Monitoring well 7-S had two detected results, which were 10.5 and 34.3 mg/L. There has been an impact on the concentration of COD in the groundwater due to the landfill.

#### Fluoride

The background concentration was calculated to be 0.193 mg/L in MW 1-S using Cohen's method. The student's t distribution was used to calculate the upper limit in MW 1-WT. This concentration was found

to be 0.757 mg/L. Of the four downgradient monitoring wells, only MW 7-S has had a significant change from the background concentrations. Monitoring well 5-S had concentrations that were less than 0.13 mg/L. The average concentration was calculated to be 0.10 mg/L in this monitoring well. The concentration of fluoride in MW 6-S ranged from 0.17 to 0.41 mg/L. The average concentration was calculated to be 0.26 mg/L. The results in MW 8-S were all below the 1-WT upper limit concentration of 0.757 mg/L. The data ranged from 0.14 to 0.26 mg/L with an average of 0.21 mg/L. There is a significant variation between the background concentrations and MW 7-S. The results ranged from 0.64 to 2.43 mg/L with an average concentration of 1.52 mg/L. The concentration of fluoride in the groundwater has been slightly impacted.

#### рH

The limits calculated for pH range from 7.16 to 8.14 in MW 1-WT and 7.22 to 7.82 in MW 1-S. In the downgradient monitoring wells, pH ranged from 6.9 to 7.8 in MW 5-S, 7.1 to 7.9 in MW 6-S, 8.9 to 10.2 in MW 7-S an 7.1 to 8.3 in MW 8-S.

#### Nitrate as N

The background concentration was calculated to be 11.07 mg/L in MW 1-S using Cohen's method. Monitoring well 5-S and 8-S had quantitative results for nitrate concentrations. In monitoring well 5-S, the three results were 1.31, 6.24, and 2.00 mg/L. Using the detection limit to calculate the average concentration for this well, the result was 1.42 mg/L. Data ranged from 1.68 to 3.01 mg/L in MW 8-S. The average concentration was calculated to be 2.48 mg/L.

A test of proportions was used to determine if there was a difference between the background MW 1-WT and the four downgradient monitoring wells, 5-S through 8-S. Nitrate was detected in three of the sample events in MW 5-S and all seven of the events in MW 8-S. The test of proportions indicated that there was no significant difference between MW 5-S and 1-WT. Monitoring well 8-S was determined to be significantly different than the background concentration.

The groundwater has been affected only slightly by the landfill with regards to nitrate concentrations.

## Sulfate

using the upper limits of the student's t 99% confidence intervals. In MW 5-S, the concentrations of sulfate ranged from 196 to 650 mg/L with an average of 385 mg/L. In MW 6-S, the average of sulfate was 235 mg/L, where the data ranged from 106 to 332 mg/L. The concentration of sulfate ranged from 336 to 494 mg/L in MW 7-S with an average concentration of 425 mg/L. The average sulfate concentration in

MW 8-S was calculated to be 803 mg/L. The results ranged from 686 to 865 mg/L. The concentration of sulfate in the groundwater has been significantly impacted by the landfill.

## Total Organic Carbon (TOC)

The background concentrations were calculated to be 4.5 and 17.1 mg/L in the MW 1-S and 1-WT using the upper limits of the student's t 99% confidence intervals. The concentration of TOC in MW 5-S ranged from below detection limits to 29 mg/L. The average concentration was calculated to be 11.7 mg/L. In MW 6-S, the data ranged from 2.01 to 13.7 mg/L. The average concentration in this well was calculated to be 10.2 mg/L. The range of TOC concentration in MW 7-S was 7 to 18 mg/L and the average was 13.4 mg/L. In MW 8-S, the results ranged from 2.05 to 14.4 mg/L with an average of 8.5 mg/L. Therefore, the landfill has had very little affect on this parameter in the groundwater.

## Total Dissolved Solids (TDS)

The background concentrations were calculated to be 540 and 488 mg/L in the MW 1-S and 1-WT, respectively, using the upper limits of the student's t 99% confidence intervals. The concentration of TDS in MW 5-S ranged from 633 to 1213 mg/L with an average concentration of 846 mg/L. In MW 6-S, results ranged from 566 to 867 mg/L with an average concentration of 797 mg/L. The range and average of TDS concentrations in MW 7-S were 802 to 1423 mg/L and 1133 mg/L, respectively. In MW 8-S, the results ranged from 1125 to 1252 mg/L. The average concentration was calculated to be 1179 mg/L for MW 8-S. The concentration of dissolved solids in the groundwater has been significantly impacted by the landfill.

## V. GROUNDWATER IMPACT- VOLATILE ORGANIC COMPOUNDS

All of the monitoring wells have been scanned for volatile organic compounds during the sampling period. Compounds have been eliminated from the list when all of the downgradient and upgradient monitoring results are below detection limit for two sampling events.

Currently there are five volatile compounds that are being analyzed for in the monitoring wells at Dixon Marquette Cement Company. These are benzene, 1,1-dichloroethane, 1,1-dichloroethene, toluene, and 1,1,1-trichloroethane. Of these compounds, benzene, 1,1-dichloroethene, toluene, and 1,1,1-trichloroethane have class I groundwater standards.

In the following sections, these five compounds are discussed with respect to the background concentrations, and the concentrations in the downgradient monitoring wells: 5-S through 8-S.

#### Benzene

The compound was not detected in the background monitoring wells MW 1-WT and 1-S. The upper background concentration limit was found to be 0.005 mg/L, the method detection limit. Benzene was not detected in the downgradient monitoring wells 5-S and 8-S.

## 1,1-Dichloroethane

The compound was not detected in the background monitoring wells MW 1-WT and 1-S. The upper background concentration limit was found to be 0.005 mg/L. In the downgradient monitoring wells 5-S through 8-S, the compound was detected in only 7-S. Seven of the eleven samples had quantitative results in MW 7-S. The average concentration was calculated to be 0.0299 mg/L.

#### 1,1-Dichloroethene

The compound was not detected in the background monitoring wells MW 1-WT and 1-S. The upper background concentration limit was found to be 0.005 mg/L, the method detection limit. In the downgradient monitoring wells, 5-S through 8-S, the compound was detected in MW 7-S only. Three of the seven samples had quantitative results. The average concentration was calculated to be 0.0058 mg/L.

#### Toluene

The sample from the shallow aquifer had 0.0082 mg/l toluene on July 24, 1998. The source of the toluene was not identified. There have been no other sampling events in whic toluene was detected in MW1-S above 0.005 mg/l. Because toluene was not confirmed in the well, its detection in the sample from July 1998 is considered to be sporadic, and toluene is not actually present in the well. Toluene was not detected in MW 1-WT, thus the upper background concentration limit is 0.005 mg/L, the method

detection limit. In the four downgradient monitoring wells, 5-S through 8-S, toluene was not detected during all of the sampling events and is not considered to be different than the background concentration.

## 1,1,1-Trichloroethane

The compound was not detected in MW 1-WT and 1-S. The upper background concentration limit was found to be 0.005 mg/L, the method detection limit. In MW 7-S, two of the sample events had results that were detected. The average concentration of 1,1,1-trichloroethane was calculated to be 0.0052 mg/L. There were no detected concentrations of 1,1,1-trichlorethane in MW 5-S, 6-S, and 8-S.

#### VI. GROUNDWATER STANDARDS

In the foll wing sections, the maximum 99.5% confidence interval of MW 5-S through 8-S is discussed with respect to the Class I Groundwater Standard. In addition, the best fit lines and trends of the sampling data are presented.

#### Aluminum

There is no Class I Groundwater Standard for this parameter. The maximum 99.5% confidence interval for MW 5-S through MW 8-S was calculated to be 2.522 mg/L.

Linear trend lines for each of the four monitoring wells indicate a decreasing slope for the data. Results from other downgradient monitoring wells indicate that the 99.5% confidence interval would be met in the future et Dixon Marquette Cement Company.

#### Antimony

The Class I Groundwater Standard for this parameter is 0.006 mg/L. The 99.5 % confidence interval was calculated to be 0.002 mg/L using data from MW 7-S.

Since the parameter was not detected in the other three wells, the trend line was steady, without a slope. The trend line for MW 7-S was negative, representing a decreasing trend. The parameter was not detected in two downgradient wells, MW 3-S, and MW 4-S.

#### Arsenic

The Class I Groundwater Standard for this parameter is 0.05 mg/L. The maximum 99.5% confidence interval was calculated to be 0.041 mg/L using data from MW 7-S.

The linear trend lines indicate a decreasing slope for each of the four monitoring wells. Although the trend lines indicate a decreasing slope for each of the monitoring wells, it is expected that the concentrations of arsenic is expected to rise to levels that exceed the standard since MW 3-S has consistently exceeded the standard.

#### Barium

The Class I Groundwater Standard for this parameter is 2.0 mg/L. The maximum 99.5% confidence interval for the four downgradient wells, 5-S through 8-S was calculated to be 0.371 mg/L.

Each of the four trend lines has a negative slope, indicating a decreasing trend in the concentrations over time. Using data from MW 3-S and 4-S and the trend lines for the four monitoring wells, the concentration of barium is not expected to rise to concentrations greater than the standard.

Beryllium

The Class I Groundwater Standard is 0.004 mg/L for beryllium. The maximum 99.5% confidence interval was calculated to be 0.002 mg/L using the data from MW 8-S.

Since the parameters were not detected in the other three wells, the trend lines were steady, without slopes. The trend line for MW 8-S was negative, representing a decreasing trend. The monitoring data from MW 3-S and 4-S indicate that there should not be an increase in the concentrations of this parameter.

Boron

The Class I Groundwater Standard is 2.0 mg/L. The maximum 99.5% confidence interval was calculated to be 1.452 mg/L from the sampling data at MW 7-S.

All of the trend lines for the four downgradient monitoring wells are negative, indicating a decreasing trend. Using the trend lines, and monitoring data from other downgradient wells, it is not expected that the concentration of boron would exceed this standard.

Cadmium

The Class I Groundwater Standard for this parameter is 0.005 mg/L. The 99.5% confidence interval was calculated to be 0.002 mg/L for the four downgradient monitoring wells.

The slopes of the trend lines indicate a that the concentrations are steady with a slight decrease in the concentration over time. Since cadmium was not detected in MW 3-S and 4-S, the landfill has had very little affect on the groundwater for this parameter.

Calcium

There is no Class I Groundwater Standard for this parameter. The maximum 99.5% confidence interval was calculated to be 420 mg/L for the four downgradient wells.

The linear trend lines have a positive slope for MW 6-S and 7-S indicating a slightly increasing trend over time. The results from MW 3-S do not indicate an increase in the parameter. Values were below 10 mg/L in MW 3-S.

Chromium-

The maximum 99.5% confidence interval was calculated to be 0.030 mg/L, this is significantly lower than the Class I Groundwater Standard of 0.1 mg/L.

The trends indicate that the concentration of chromium slightly is decreasing in MW 5-S through 7-S. There is a small increasing trend in MW 8-S. There is very little difference between the concentrations of MW 5-S through 8-S and the chromium concentrations in MW 3-S.

#### Cobalt

The 99.5% confidence interval for the four downgradient wells was calculated to be 0.027 mg/L. This is significantly lower than the Class I Groundwater Standard of 1 mg/L.

All of the trend lines for the four monitoring wells indicate that the cobalt concentrations are remaining steady with a slight decrease over time. The data from MW 3-S indicate that the landfill is having very little affect on this parameter, and one would not expect a significant increase over time.

## Copper

The 99.5% confidence interval for the four monitoring wells 5-S through 8-S was calculated to be 0.084 mg/L. This is significantly lower the Class I Groundwater Standard of 0.65 mg/L.

Except for MW 6-S, the trends indicate a slightly decreasing slope. The slope for MW 6-S indicates an increasing trend over time. This, however, is a very shallow slope.

#### Lead

The maximum 99.5% confidence interval was calculated to be 0.012 using data from MW 6-S. This is approximately 60% higher than the class I groundwater standard.

The linear trends for the four monitoring wells have negative slopes, indicating a decreasing trend. Concentrations of lead in MW 3-S were below 0.002 mg/L, which indicate that the landfill has not had a significant affect on this parameter.

## Iron

The maximum 99.5% confidence interval for the four downgradient wells was calculated to be 6.399 mg/L using the data from MW 6-S. This is approximately 128% increase from the Class I Groundwater Standard of 5 mg/L.

The four best-fit lines have negative slopes, representing a decreasing trend. Concentrations of iron in MW 3-S are less than 1 mg/L, whereas the concentrations of MW 4-S range from 0.323 to 5.39 mg/L,

#### Magnesium

There is no Class I Groundwater Standard for this parameter. The maximum 99.5 % confidence interval was calculated to be 171.5 mg/L using data from MW 5-S.

The best fit line for the sampling data in MW 5-S, 7-S and 8-S indicate an increasing trend for this parameter. Monitoring well 3-S has concentrations that are significantly less than the 1-WT background concentration.

Manganese

The Class I Groundwater Standard is 0.15 mg/L for this parameter. The maximum 99.5% confidence interval was calculated to be 3.251 mg/L using data from MW 6-S.

Only MW 8-S was represented by a positive slope during trend analysis. This indicates that the concentration manganese is increasing in this monitoring well.

Mercury

The class I ground water standard for this parameter is 0.002 mg/L. Since the parameter was not detected, the lowest non-detectable limit, 0.0002 mg/L was found to be the maximum 99.5% confidence interval.

Nickel

The 99.5% confidence interval was calculated to be 0.044 mg/L, this is significantly lower than 0.1, the Class I Groundwater Standard. The best fit lines four each of the four monitoring wells indicate a slightly decreasing trend in the data.

Potassium

There is no class I ground water standard for this parameter. The maximum 99.5% confidence interval was calculated to be 838.1 mg/L.

Although the four best-fit lines that represent the data have negative slopes, or decreasing trends, it is unlikely that the concentration in these wells will decrease. Data from MW 3-S suggest that the concentration will increase as the average concentration in this well is 1636 mg/L.

Selenium

The maximum 99.5% confidence interval was calculated to be 0.011 mg/L. There is approximately 355% difference between the confidence interval and the Class I Groundwater Standard of 0.05 mg/L.

Three of the four best-fit lines have a positive slope, indicating an increasing trend. The trend line for MW 5-S is negative. Data from MW 3-S indicate that there may be an increase in the concentrations of selenium over time.

Sodium

There is no Class I Groundwater Standard for this parameter. The maximum 99.5% confidence interval was calculated to be 74.4 mg/L using the data from MW 5-S.

Three of the four trend lines indicate that the concentration of sodium is increasing over time. Data from MW 3-S further indicate that sodium concentrations are being affected by the landfill and the concentration will increase in the four downgradient monitoring wells. The average concentration in MW 3-S is 122.1 mg/L.

#### Thallium

The Class I Groundwater Standard for thallium is 0.002 mg/L. The 99.5% confidence interval was calculated to be 0.004 mg/L using data from MW 8-S.

There were no trend lines for the monitoring wells 5-S through 7-S. The best-fit line for MW 8-S has a positive slope, indicating an increasing trend. Data from MW 3-S indicate that an increase in thallium is unlikely to occur.

#### Vanadium

There is no Class I Groundwater Standard for this parameter. The maximum 99.5% confidence interval was calculated to be 0.021 mg/L using the data from MW 7-S.

Three of the four slopes of the best-fit lines were negative, indicating a decreasing trend. The slope on the trend line for MW 8-S was positive. Data from MW 3-S indicate that there is a slight increase in vanadium concentrations over time, as the average concentration in this well is 0.012 mg/L.

#### Zinc

The 99.5% confidence interval was calculated to be 0.268 mg/L, this is significantly lower than the groundwater standard of 5 mg/L.

The trend lines for the four monitoring wells indicate a decrease in concentration of zinc over time. Data from MW 3-S indicate that the landfill has had little affect on the zinc concentrations, as the average is 0.011 mg/L.

#### Total Alkalinity

There is no Class I Groundwater Standard for this parameter. The 99.5% confidence interval was calculated to be 578 mg/L, using data from MW 5-S.

The best fit lines for the four downgradient wells, 5-S through 8-S, were positive, indicating an increase in the concentration of total alkalinity over time. Data from MW 3-S indicates that the alkalinity in these wells will continue to rise as the average concentration in this well is 1488 mg/L.

### Phenol Alkalinity

There is no Class I Groundwater Standard for this parameter. The 99.5% confidence interval was calculated to be 305 mg/L using data from MW 7-S.

The slope of the best-fit line for the results in MW 7-S indicates a positive increase in concentration of phenol alkalinity over time. In addition, MW 3-S has concentrations that average 765 mg/L.

#### Bicarbonate Alkalinity

There is no Class I Groundwater Standard for this parameter. The 99.5% confidence interval was calculated to be 528 mg/L using data from MW 5-S.

The best-fit lines for MW 5-S and 6-S have positive slopes, indicating an increase in concentration over time. Data from MW 3-S also indicates that there will be a steady increase in the bicarbonate alkalinity as the range of sampling data is 120 to 1450 mg/L and the average concentration is 523 mg/L.

## Ammonia Nitrogen

There is no Class I Groundwater Standard for this parameter. The 99.5% confidence interval was calculated to be 0.70 mg/L, using data from MW 7-S.

The trends for MW 5-S and MW 7-S indicate an increase in ammonia nitrogen over time. Data from MW 3-S had two sampling events with detectable concentrations. The results were 1.60 and 0.61 mg/L.

#### Chloride

The Class I Groundwater Standard is 200 mg/L, whereas the 99.5% confidence interval was calculated to be 241 mg/L for the four downgradient monitoring wells, 5-S through 8-S. This is a 17% difference.

Only the best fit line for the *iAW* 5-S sampling data indicates a decreasing trend over time. Chloride concentrations are expected to continue to rise in the four downgradient monitoring wells. Monitoring well 3-S has chloride concentrations that range from 130 to 231 mg/L, the average was calculated to be 169.0 mg/L.

### Chemical Oxygen Demand (COD)

There is no Class I Groundwater Standard for COD. The maximum 99.5% confidence interval was calculated to be 155.3 mg/L using data from MW 5-S.

All four best-fit lines indicated that the concentration of COD would increase over time. Monitoring well 3-S also indicates that the concentration of COD will increase in the four downgradient monitoring wells. The concentration of COD ranged from below detection limits to 242 mg/L.

#### Fluoride

The Class I Groundwater Standard for fluoride is 4 mg/L. The maximum 99.5% confidence interval was calculated to be 2.48 mg/L using data from MW 7-S.

The best fit lines indicate that the concentration of fluoride is increasing in all of the monitoring wells except for MW 7-S. In addition, the concentration of fluoride in MW 3-S ranged from 3.02 to 4.44 mg/L with an average concentration of 3.60 mg/L. This further indicates that there will be an increase in the fluoride concentration in the four downgradient monitoring wells.

#### Nitrate as N

The Class I Groundwater Standard for nitrate is 10 mg/L. The 99.5% confidence interval was calculated to be 4.84 mg/L. There is approximately 107% difference between the two values.

The trend lines showed a decreasing slope for MW 5-S and an increasing one for MW 8-S. Since the concentrations in MW 3-S are similar to the MW 5-S, the concentration of nitrate is not expected to increase.

#### pH

The two sided 99% confidence interval was calculated to be 7.0 to 10.4 for the four downgradient monitoring wells 5-S through 8-S. Although the linear trends indicate a decreasing pH, values greater than 10 were reported in MW 7-S.

#### Sulfate

The Class I Groundwater Standard for sulfate is 400 mg/L. The 99.5% confidence interval was calculated to be 897 mg/L from the MW 8-S data.

The best-fit lines indicate that sulfate is increasing in MW 8-S only. Results from MW 3-S indicate that the concentration of sulfate will increase. The average concentration was calculated to be 1076 mg/L in MW 3-S.

## Total Organic Carbon (TOC)

There is no Class I Groundwater Standard for total organic carbon. The maximum 99.5% confidence interval was calculated to be 28.9 mg/L.

All of the best-fit lines indicate a decreasing trend in concentration over time. Data from MW 3-S indicates that the concentration of TOC will increase as the average concentration is 17.32 mg/L.

## Total Dissolved Solids (TDS)

The Class I Groundwater Standard for this parameter is 1200 mg/L. The maximum 99.5% confidence interval for the four downgradient monitoring wells was calculated using results from MW 7-S. The result is 1478 mg/L.

The best-fit line for MW 8-S indicates an increasing trend for the concentration of TDS over time. The other three have a negative slope, indicating a decrease in concentration. Data from MW 3-S indicates that the concentration of TDS will increase instead of decrease. The average concentration is 3843 mg/L.

## Benzene

The class I groundwater standard for this compound is 0.005 mg/L. The maximum 99.5% confidence interval could not be calculated for monitoring wells 5-S through 8-S since benzene was not detected during the sampling events. All of the trend lines are represented by horizontal lines that have no slope.

#### 1,1-Dichloroethane

There is no class I groundwater standard for this parameter. The maximum 99.5% confidence interval was calculated to be 0.0669 mg/L using data from MW 7-S. Since the compound was not detected in MW 5-S, 6-S, and 8-S, the trend lines are horizontal without a slope. The best fit line in MW 7-S was represented by a positive slope representing an increasing trend.

#### 1,1-Dichloroethene

The class I groundwater standard for this parameter is 0.007 mg/L. The maximum 99.5% confidence interval was calculated to be 0.0076 mg/L using data from MW 7-S. Since the 1,1-dichloroethene was not detected in MW 5-S, 6-S, and 8-S, the trend lines are horizontal without a slope. The best fit line in MW 7-S was represented by a positive slope representing an increasing trend.

## Toluene

The class I groundwater standard for this compound is 1 mg/L. The maximum 99.5% confidence interval could not be calculated for monitoring wells 5-S through 8-S since toluene was not detected during the sampling events. All of the trend lines are represented by horizontal lines that have no slope.

#### 1,1,1-Trichloroethane

The class I groundwater standard for this compound is 0.2 mg/L. The maximum 99.5% confidence interval was calculated to be 0.0059 mg/L using data from MW 7-S. Since the 1,1,1-trichloroethane was not detected in MW 5-S, 6-S, and 8-S, the trend lines are horizontal without a slope. The best fit line in MW 7-S was represented by a negative slope representing a decreasing trend.

# VII. ADJUSTED GROUNDWATER STANDARDS

In the following table, the Class Groundwater Standards, the maximum 99.5% Confidence Interval of MW 5-S through 8-S, the upper limits of the 99<sup>th</sup> percentile for MW 1-WT and 1-S, and the calculated 99.5% confidence interval for MW 3-S are given. The requested adjusted groundwater standard is listed. All concentrations are given in mg/L, unless noted otherwise.

	Class I Standard	MW 5S - 8S Max 99.5% CI	MW 1-S Upper Limit	MW 1-WT Upper limit	MW 3-S 99.5% CI	Requested Adjusted Standard
Aluminum		<u>2.522</u>	0.04	0.454	0.730	2.522
Antimony	0.006	0.002	Tst of Prop	0.001	0.001	0.006
Arsenic	0.05	0.041	Tst of Prop	· 0.003	0.141	0.141
Barium	2	0.371	0.091	0.227	0.068	2
Beryllium	0.004	0.002	0.001	0.001	0.001	0.004
Boron	2	1.452	Tst of Prop	Tst of Prop	0.212	2
Cadmium	0.005	0.002	0.001	0.001	0.006	0.005
Calcium		420	109	126	6.6	420
Chromium	<u>0.1</u>	0.030	0.002	0.002	0.032	0.1
Cobalt	1	0.027	0.003	0.008	0.003	1
Copper	0.65	0.084	0.005	0.012	0.037	0.65
Lead	0.0075	0.012	0.001	0.006	0.004	0.012
Iron	5	<u>6.399</u>	0.465	0.891	1.017	6.4
Magnesium		<u>171.5</u>	53.9	50.4	4.6	171.5
Manganese	0.15	3.251	0.045	2.687	0.186	3.3
Mercury	0.002	0.0002	0.0002	Tst of Prop	0.0011	0.002
Nickel	<u>0.1</u>	0.044	0.007	0.023	0.031	0.1
Potassium		838	Tst of Prop	4.004	2309	2309
Selenium	0.05	0.011	0.004	Tst of Prop	0.012	0.05
Sodium		74.4	12.8	53.7	140.5	140.5
Thallium	0.002	0.004	Tst of Prop	Tst of Prop	0.002	0.004
Vanadium		0.021	Tst of Prop	0.007	0.026	0.02
Zinc	<u>5</u>	0.268	0.058	0.014	0.018	5
Alkalinity, total		578	253	342	1699	1699
Alkalinity, phenol.		305.0	Tst of Prop	1.0	1121	1121
Alkalinity, bicarb		528	250.036	334.029	977	977
Ammonia nitrogen		0.70	Tst of Prop	Tst of Prop	1.25	1.25
Chloride	200	241	30.8	25.4	197	241
COD		155	5.0	Tst of Prop	187	187
Fluoride	4	2.48	0.19	0.76	4.42	4,4
Nitrate as N	10	4.84	11.07	Tst of Prop	0.88	11,1
pH (stnd units)		7.0 to 10.4	7.22 to 7.82	7.16 to 8.14	9.0 to 10.9	7.0 to 10.9
Sulfate	400	897	178	243	1245	1245
TOC		28.9	4.5	17.1	26.51	29
Ttl Dis Solids	1200	1478	540	488	4338	4338

	Class I Standard	MW 5S - 8S Max 99.5% CI	MW 1-S Upper Limit	MW 1-WT Upper limit	MW 3-S 99.5% CI	Requested Adjusted Standard
Benzene	<u>0.005</u>	NA	0.005	0.605	NA	0.065
1,1- dichloroethane		0.0669	0.005	0.005	5.4 mg/L	0.7 *
1,1- dichloroethene	0.007	0.0076	0.005	0.005	NA	0.035 **
Toluene	L	NΛ	0.005	0.005	NA	1
i,1,1- trichloroethane	0.2	0.0059	0.005	0.005	12.0	0.2

NA- Concentrations for all sampling events were below detection limits. The 99.5% confidence interval was not calculated for this parameter.

<sup>\*</sup> The Class I groundwater remediation objective from Section 742 was used because a groundwater standard is not available.

<sup>\*\*</sup> The Class II 620 groundwater standard is proposed.

Due to the concentrations in monitoring well 3-S, some of the parameters are expected to increase above the maximum confidence interval of 5-S through 8-S. These parameters include: arsenic, potassium, sodium, alkalinity, sulfate, and total dissolved solids. In these instances, the 99.5% confidence interval of 3-S was requested as the adjusted standard. In the following paragraphs, each of these parameters is explained with respect to the concentrations in MW 3-S and the maximum 99.5% confidence interval of MW 5-S through 8-S.

Arsenic—The class I groundwater standard is 0.05 mg/L, which is significantly higher than the 99.5% confidence interval calculated for monitoring wells 5-S through 8-S. Seven of the nine sample results in MW 3-S were greater than the standard of 0.05 mg/L. The average concentration was calculated to be 0.083 mg/L.

Potassium— The maximum 99.5% confidence interval of MW 5-S through 8-S is 838 mg/L. Potassium concentrations in MW 3-S ranged from 200 to 2314 mg/L with an average concentration of 2010 mg/L.

Sodium—The maximum 99.5% confidence interval of MW 5-S through 8-S is 74.4 mg/L. This concentration is significantly lower than the average concentration of sodium in MW 3-S. Sodium ranged from 82 to 145 mg/L with an average of 122.1 mg/L in this downgradient well.

Total Alkalinity—The maximum 99.5% confidence interval of MW 5-S through 8-S is 578 mg/L. In monitoring well 3-S, total alkalinity peaked at 1660 mg/L on November 29, 2000. The concentration of total alkalinity has been consistently rising in this well. The average was calculated to be 1488 mg/L.

Phenol Alkalinity— The average concentration in MW 3-S was calculated to be 765 mg/L, however the peak concentration of phenol alkalinity was measured to be 1155 mg/L during the June 2000 sampling event. The maximum 99.5% confidence interval of MW 5-S through 8-S is 305 mg/L.

Bicarbonate Alkalinity—The average concentration of bicarbonate alkalinity in MW 3-S is 523 mg/L, this is lower than the maximum 99.5% confidence interval of MW 5-S through 8-S, 528 mg/L. Since the maximum concentration in MW 3-S is 1450 mg/L, the four downgradient monitoring wells are expected to exceed the adjusted groundwater standard.

Sulfate—The maximum 99.5% confidence interval of MW 5-S through 8-S is 897 mg/L. This is significantly lower then the average concentration in MW 3-S. The average concentration was calculated to be 1076 mg/L and the peak concentration in this well was 1400 mg/L.

Total Dissolved Solids – The maximum 99.5% confidence interval is 1478 mg/L in MW 5-S through 8-S. This concentration is significantly lower than the average concentration, 3843 mg/L, in MW 3-S. Concentrations of total dissolved solids have been increasing in MW 3-S over time. The peak concentration was 4400 mg/L during the September 2000 sampling event.

Since there were exceedances of the requested adjusted groundwater standard in MW 3-S, the following parameters are also considered to be of future concern.

Vanadium – During the sampling period, there was one MW 3-S concentration of vanadium that was greater than the 99.5% confidence interval of 0.021 mg/L. This occurred during the November 2000 sampling event and the concentration was 0.029 mg/L.

Ammonia Nitrogen – Two of the seven sample results had concentrations of ammonia nitrogen in MW 3-S. The concentrations were 1.60 and 0.61 mg/L. The requested adjusted standard for this parameter is 0.70 mg/L. Thus one sample was significantly greater than the standard.

Chemical Oxygen Demand-Although the average concentration of COD in MW 3-S is less than the requested standard, there is some concern with future results. During the June 2000 sampling event, the concentration of COD peaked at 242 mg/L.

Fluoride- In monitoring well 3-S, one concentration was greater than the Class I Groundwater Standard. The concentration was 4.44 mg/L, a 110% increase above the standard.

# **SUMMARY TABLES**

Parameter	MW 5-S	MW 6-S	MW 7-S	MW 8-S	Class I GroundWater Stnd	Maximum Value
Aluminum	0.277	2.522	0.844	0.664	Ground Water Stud	2,522
Antimony	0.001	0.001	0.002	0.004	0.006	0.002
Arsenic	0.004	0.004	0.041	0.004	0.05	0.041
Barium	0.269	0.371	0.057	0.141	2	0.371
Beryllium	0.001	0.001	0.001	0.002	0.004	0.002
Boron	0.786	0.751	1.452	0.318	2	1.452
Cadmium	0.002	0.001	0.001	0.002	0.005	0.002
Calcium	318	420	45	268		420
Chromium	0.006	0.006	0.005	0.030	0.1	0.030
Cobalt	0.005	0.027	0.001	0.025	1	0.027
Copper	0.014	0.027	0.009	0.084	0.65	0.084
Lead	0.001	0.012	0.003	0.004	0.0075	0.012
Iron	1.834	6.399	1.806	4.186	5	6.399
Magnesium	171.5	95.8	15.9	119.3		171.5
Manganese	0.913	3.251	0.187	3.133	0.15	3,251
Mercury	0.0002	0.0002	0.0002	0.0002	0.002	0.0002
Nickel	0.016	0.039	0.016	0.044	0.1	0.044
Potassium	259. <b>6</b>	<b>77.</b> 6	838.1	302		838
Selenium	0.005	0.005	0.004	0.011	0.05	0.011
Sodium	74.4	68.0	53.5	62.2		74.4
Thallium	0.001	0.001	0.001	0.004	0.002	0.004
Vanadium	0.018	0.019	0.021	0.004		0.021
Zinc	0.039	0.035	0.033	0.268	5	0.268
Alkalinity, total	578	425	534	317	•	578
Alkalinity, phenol.	1	2.4	305	1.0		305.0
Alkalinity, bicarb	528	409	441.57	309		528
Ammonia nitrogen	0.16	0.10	0.7	0.10		0.70
Chloride	132.0	241	119.32	206	200	241
COD	155.3	127	31.24	119		155
Fluoride	0.15	0.39	2.48	0.28	4	2.48
Nitrate as N	4.84	0.10	0.10	3.27	10	4.84
pH Upper Limit (stnd units)	7.9	8.1	10.4	8.3		10.4
pH Lower Limit (stnd units)	7.0	7.0	9.0	7.0		7.0
Sulfate	694	352	523	897	400	897
TOC	28.89	20.83	19.3	16.5		28.9
Ttl Dis Solids	1225	957	1478	1245	. 1200	1478
Benzene	NA	NA	NA	NA	0.005	NA
1,1-dichloroethane	NA	NA	0.0669	NA		0.0669
1,1-dichloroethene	NA	NA	0.0076	NA	0.007	0.0076
Toluene	NA	NA	NA	NA	1	NA
1,1,1-trichloroethane	NA	NA	0.0059	NA	0.2	0.0059
	All values	in mg/L un	less noted o	therwise.		

· Parameter	MW 3-S	MW 5-S	MW 6-S	MW 7-S	MW 8-S		Class I GroundWater Stnd	Maximum Value
Aluminum	0.730	0.277	2.522	0.844	0.664		÷ .	2.522
Antimony	0.001	0.001	0.001	0.002	0.001		0.006	0.002
Arsenic	0.141	0.004	0.004	0.041	0.004		0.05	0.141
Barium	0.068	0.269	0.371	0.057	0.141		2	0.371
Beryllium	0.001	0.001	0.001	0.001	0.002		0.004	0.002
Boron	0.212	0.786	0.751	1.452	0.318		2	1.452
Cadmium	0.006	0.002	0.001	0.001	0.002		0.005	0.006
Calcium	6.6	318	420	45	268			420
Chromium	0.032	0.006	0.006	0.005	0.030		0.1	0.032
Cobalt	0.003	0.005	0.027	0.001	0.025	**	1	0.027
Copper	0.037	0.014	0.027	0.009	0.084	*	0.65	0.084
Lead	0.004	0.001	0.012	0.003	0.004		0.0075	0.012
Iron	1.017	1.834	6.399	1.806	4.186		5	6.399
Magnesium	4.6	171.5	· 95.8	15.9	119.3			171.5
Manganese	0.186	0.913	3.251	0.187	3.133		0.15	3.251
Mercury	0.0011	0.0002	0.0002	0.0002	0.0002		0.602	0.0011
Nickel	0.031	0.016	0.039	0.016	0.044		0.1	0.044
Potassium	2309	259.6	77.6	838.1	302			2309
Selenium	0.012	0.005	0.005	0.004	0.011		0.05	0.012
Sodium	140.5	74.4	68.0	53.5	62.2			140.5
Thailium	0.002	0.001	0.001	0.001	0.004		0.002	0.004
Vanadium	0.026	0.018	0.019	0.021	0,004			0.026
Zinc	0.018	0.039	0.035	0.033	0.268		5	0.268
								0.000
								0.00
Alkalinity, total	1699	578	425	534	317			1699
Alkalinity, phenol.	1120.8	1	2.4	305	1.0			1121
Alkalinity, bicarb	977	528	409	441.57	309			977
Ammonia nitrogen	1.25	0.16	0.10	0.7	0.10			1.25
Chloride	197.5	132.0	241	119.32	206		200	241
COD	187.4	155.3	127	31.24	119			187
Fluoride	4.42	0.15	0.39	2.48	0.28		4	4.42
Nitrate as N	0.88	4.84	0.10	0.10	3.27		10	4.84
pH Upper Limit (stnd units)	10.9	7.9	8.1	10.4	8.3			10.9
pH Lower Limit (stnd units)	10.3	7.0	7.0	9.0	7.0			7.0
Sulfate	1245	694	· 352	523	897		400	1245
TOC	26.51	28.89	20.83	19.3	16.5			28.9
Ttl Dis Solids	4338	1225	957	1478	1245		1200	4338
Benzene	NA	NA	NA	NA	NA		5	NA
1,1-dichloroethane	0.005	NA	NA	0.0669	NA			0.0669
1,1-dichloroethene	NA	NA	NA	0.0076	NA		7	0.0076
Toluene	NA	NA	NA	NA	NA		1000	NA
1,1,1-trichloroethane	0.012	NA	NA	0.0059	NA		200	0.012
•	All values	in mg/L un	less noted o	otherwise.				

	Class I .Grndwtr Standard	MW 5S to 8S Max 99.5% CI	MW 1-S Upper Limit	MW 1-WT Upper limit	MW 3-8 99.5% CI	Requested Adjusted Groundwater Stnd
Aluminum		2.522	0.04	0.454	0.730	2.522
Antimony	0.006	0.002	Tot of Prop	0.001	0.001	0.006
Arsenic	0.05	0.041	Tst of Prop	0.003	0.141	0.141
Barium	2	0.371	0.091	0.227	0.068	2
Beryllium	0.004	0.002	0.001	0.001	0.001	0.004
Boron	2	1.452	Tst of Prop	Tst of Prop	0.212	2.000
Cadmium	0.005	0.002	0.001	<b>0</b> .001	0.006	0.005
Calcium	-	420	109 .	126	6.6	420
Chromium	0.1	0.030	0.002	0.002	0.032	0.1
Cobalt	. 1	0.027	0.003	0.008	0.003	1
Copper	0.65	0.084	0.005	0.012	0.037	0.65
Lead	0.0075	0.012	0.001	0.006	0.004	0.012
Iron	5	6.399	0.465	0.891	1.017	6.399
Magnesium		171.5	53.903	50.478	4.6	171.5
Manganese	0.15	3.251	0.045	2.687	0.186	3.251
Mercury	0.002	0,0002	0.0002	Tst of Prop	0.0011	0.002
Nickel	0.1	0.044	0.007	0.023	0.031	0.1
Potassium	•••	838	Tst of Prop	4.004	2309	2309
Selenium	0.05	0.011	0.004	Tst of Prop	0.012	0.05
Sodium	0.00	74.4	12.8	53.7	140.5	140.5
Thallium	0.002	0.004	Tst of Prop	Tst of Prop	0.002	0.004
Vanadium	0.002	0.021	Tst of Prop	0.007	0.026	0.026
• Zinc	5	0.268	0.058	0.014	0.018	5
Alkalinity, total		578	253	342	1699	1699
Alkalinity, phenol.		305.0	Tst of Prop	1.0	1121	1121
Alkalinity, bicarb		528	250.036	334.029	977	977
Ammonia nitrogen		0.70	Tst of Prop	Tst of Prop	1.25	1.25
Chloride	200	241	30.8	25.4	197	241
COD		155	5.0	Tst of Prop	187.4	187
Fluoride	4	2.48	0.19	0.76	4.42	4.42
Nitrate as N	10	4.84	11.07	Tst of Prop	0.88	11.07
pH Upper Limit (stud units)		10.4	7.82	8.14	10.9	10.9
pH Lower Limit (stnd units)		7.0	7.22	7.16	10.3	7.0
Sulfate	400	897	178	243	1245	1245
тос		28.9	4.5	17.1	26.51	28.9
Ttl Dis Solids	1200	1478	540	488	4338	4338
Benzene	0.005	NA	0.005	0.005	NA	0
1,1-dichloroethane		0.0669	0.905	0.005	0.005	0.7*
1,1-dichloroethene		0.0076	0.005	0.005	NA	0.35**
Toluene	1	NA	0.005	0.005	NA	1
1,1,1-trichloroethane	0.2	0.0059	0.005	0.005	0.012	0.2
	All values in m	g/L unless noted oil	erwise.			

<sup>\*</sup> The Class I groundwater remediation objective from Section 742 was used because a groundwater standard is not available.

<sup>\*\*</sup> The Class II 620 groundwater standard is proposed.

• •	MW 3-S AVE	MW 5-S AVE	MW 6-S AVE	. MW 7-S AVE	MW 8-S AVE	Class I Gradwtr Standard	MW 5S to 8S Max 99.5% Cl	MW 1-S Upper Limit	MW 1-WT Upper limit
Aluminum	0.322	0.168	0.754	0.387	€.326		2.522	0.04	0.454
Antimony	0.001	0.001	0.001	0.001	0.001	0.006	0.002	Tst of Prop	0.001
Arsenic	0.083	0.002	0.002	0.023	0.002	0.05	0.041	Tst of Prop	0.003
Barium	0.051	0.138	0.258	0.042	0.075	2	0.371	0.091	0.227
Beryllluan	0.001	0.001	0.001	0.001	0.001	0.004	0.002	0.001	0.001
Boron	0.107	0.464	0.424	0.436	0.208	2	1.452	Tst of Prop	Tst of Prop
Cadmium	0.002	0.001	0.001	0.001	0.001	0.005	0.002	0.001	0.001
Calcium	4.3	240	234	25	237		420	109	126
Chromium	0.012	0.003	0.003	0.003	0.010	0.1	0.030	0.002	0.002
Cobalt	0.001	0.002	0.012	0.001	0.006	1	0.027	0.003	0.008
Copper	0.017	0.006	0.013	0.005	0.024	0.65	0.084	0.005	0.012
Lead	0.002	0.001	0.004	0.002	0.003	0.0075	0.012	0.001	0.006
Iron	0.576	0.898	2.665	0.805	1.654	5	6.399	0.465	0.891
Magnesium	2.6	82.3	55.2	9.3	100.1		171.5	53.9	50.5
Manganese	0.096	0.229	1.695	0.094	0.763	0.15	3.251	0.045	2.687
Mercury	0.0004	0.0002	0.0002	0.0002	0.0002	0.002	0.0002	0.0002	Tst of Prop
Nickel	0.016	0.005	0.020	0.010	0.014	0.1	0.044	0.007	0.023
Potassium	1636	86.0	54.8	632.6	245		838	Tst of Prop	4.0
Selenium	0.007	0.002	0.003	0.002	0.005	0.05	0.011	0.004	Tst of Prop
Sodium	122.1	37.9	46.4	40.9	39.7		74.4	12.8	53.7
Thallium	0.001	0.001	0.001	0.001	0.002	0.002	0.004	Tst of Prop	Tst of Prop
Vanadlum	0.012	0.007	0.008	0.013	0.003		0.021	Tst of Prop	0.007
Zinc	0.011	0.017	0.015	0.016	0.075	5	0.268	0.058	0.014
Alkalinity, total	1488	442	389	422	287		578	253	342
Alkalinity, phenol.	765.2	1.0	0.8	74	1.0		305.0	Tst of Prop	1.0
Alkalinity, bicarb	523	410	366	294.71	276		528	250	334
Ammonia nitrogen	0.39	0.11	0.10	0.2	0.10		0.70	Tst of Prop	Tst of Prop
Chloride	169.0	<b>62.7</b>	153	82.59	153	200	241	31	25
COD	68.3	53.2	57	20.69	41		155	5.0	Tst of Prop
Fluoride	3.60	0.10	0.26	1.52	0.21	4	2.48	0.19	0.76
Nitrate as N	0.43	1.42	0.10	0.10	2.48	10	4.84	11.07	Tst of Prop
pH Upper Limit (stnd units)	10.6	7.4	7.6	9.7	7.6		10.4	7.32	*8.14
pli Lower Limit (stnd units)	10.6	7.4	7.6	9.7	7.6		7.0	7.22	7.16
Sulfate	1076	385	235	425	803	400	897	178	243
тос	17.32	11.68	10.19	13.4	8.5		28.9	4.5	17.1
Ttl Dis Solids	3843	846	797	1133	1179	1200	1478	540	488
Benzene	NA	NA	NA	NΛ	NA	0.005	NΛ	0.005	0.005
1,1-dichloroethane	0.0045	NΛ	NA	0.0299	NA		0.0669	0.005	0.005
1,1-dichloroethene	NΛ	NΛ	NΛ	0.0058	NA	0.007	0.0075	0.005	0.005
Toluene	NA	NΛ	NΛ	NΛ	NA	1	NΛ	0.005	0.005
1,1,1-trichioroethane	0.0081	NA "	NA	0.0052	NΑ	0.2	0.0059	0.005	0.005
	All values in	mg/L unless	noted other	wise.					

# **BACKGROUND TABLES**

-	Sample 1	Sample 2	Sample 3	Sample 4	Sample 5	Sample 6
Aluminum	<0.1	<0.1	<0.04	<0.04	<0.04	< 0.04
Antimony	<0.1	< 0.1	< 0.001	< 0.001	<0.001	<0.001
Arsenic	< 0.005	< 0.08	< 0.08	<0.08	<0.001	< 0.001
Barium	0.067	0.064	0.056	0.069	0.072	0.067
Beryllium	< 0.01	< 0.01	<0.01	< 0.001	< 0.001	<0.001
Boron	< 0.1	<0.05	< 0.05	<0.05	0.05	< 0.05
Cadmium	< 0.005	< 0.02	< 0.02	< 0.02	< 0.001	<0.001
Calcium	95	91	100	100	117	98.3
Chromium	0.001	0.001	0.001	0.001	0.001	0.003
Cobalt	0.001	0.001	0.001	0.004	0.001	0.061
Copper	0.002	0.002	0.002	0.005	0.001	0.004
Lead	< 0.005	<0.1	<0.1	<0.1	<0.001	<0.001
Iron	0.258	0.298	0.153	0.2	0.201	0.22
Magnesium	48	45	49	59.7	48	50.7
Manganese	0.031	0.029	0.025	0.03	0.029	0.045
Mercury	< 0.0002	< 0.0002	< 0.0002	<0.0002	< 0.0002	< 0.0002
Nickel	0.005	0.006	0.006	<0.058	. <0.05	<0.05
Potassium	1.04	0.89	1.81	1.8	1.13	2.2
Sclenium	0.004	0.003	0.003	< 0.005	< 0.15	< 0.15
Sodium	11	10	11	11	12.5	11.3
Thallium		<1	<1	< 0.001	< 0.001	<0.001
Vanadium	< 0.05	< 0.05	0.001	< 0.001	< 0.001	<0.001
Zine	0.08	0.039	0.034	0.032	0.037	0.01
Alkalinity, total	256	240	250	188	216	242
Alkalinity, phenol.		<5	<5	4	<1.0	<1
Alkalinity, bicarbonate	256	240	250	170	201	230
Ammonia nitrogen	<.2	<.2	<.2	<.1	0.27	<0.1
Chloride	31	30	31	31	32	29.6
COD		<5	<5	<5	<20	<20
Fluoride	0.18	0.15	0.16	<0.2	<0.2	<0.21
Nitrate	6.9	6.9	9.7	7.5	6.03	6.89
pH Upper Limit	7.41	7.3	7.1	7.4	7.4	7.8
pH Lower Limit						1/0
Sulfate	160	180	150	180	170	169
тос	3.9	1	5	6.09	2.32	1.45
TDS	574	550	580	550	550	428
Benzene	<0.0010	<0.0005	<0.0005	<0.0005	<0.0050	<0.0050
1,1-dichloroethane	< 0.0010	< 0.0010	<0.0010	<0.0010	<0.0050	<0.0050
1,1-dichloroethene	<0.0010	< 0.0020	< 0.0020	<0.0020	<0.0050	<0.0050
Toluene	<0.0010	0.0082	<0.0010		<0.0050	<0.0050
1,1,1-trichloroethane	<0.0010	< 0.0010	<0.0010	<0.0010	<0.0050	<0.0050
•	All concen	trations in mg	/L.			

	Sample 7	Sample 8	Sample 9	Sample 10	Sample 11	Sample 12	Sample 13
Aluminum	< 0.04	< 0.04	< 0.04				•
Antimony	< 0.001	0.001	< 0.001				
Arsenic	< 0.001	< 0.001	< 0.001	0.001	< 0.001		
Barium	0.066	0.075	0.062	0.069	0.135		
Beryllium	< 0.001	< 0.001	< 0.001	< 0.001			
Boron	< 0.05	0.078	< 0.05	0.048			
Cadmium	< 0.001	< 0.001	< 0.001	< 0.001	< 0.001		
Calcium	103	112	100	106	104		
Chromium	<0.04	< 0.02	< 0.02	< 0.02	< 0.001		
Cobalt	< 0.02	<0.02	< 0.02	< 0.001			
Copper	< 0.02	< 0.02	< 0.02	< 0.001			
Lead	< 0.001	< 0.001	< 0.001	< 0.001	< 0.001		
lron	1.01	< 0.05	< 0.1	< 0.1	<0.1	< 0.1	
Magnesium	52.3	49.5	51.5	48,5			
Manganese	0.039	0.044	0.028	0.032	0.063	0.0005	
Mercury	< 0.0002	< 0.0002	< 0.0002	< 0.0002			
Nickel	0.003	0.004	0.005	0.01			
Potassium	<2	<2	<1	<2	<2	< 0.05	
Selenium	<0.15	0.002	<0.001	0.007	0.012		
Sodium	11.8	12.6	14.5	12.5	11.1		
Thallium	< 0.001	< 0.001	0.005	<0.001			
Vanadium	0.002	0.002	0.008				
Zinc	0.019	0.007	0.036	0.067			
Alkalinity, total	240	240	230	246			
Alkalinity, phenol.	<1						
Alkalinity, bicarbonate	228	227	226	238			
Ammonia nitrogen	< 0.10	<0.1	0.23	<0.1			
Chloride	27.5	27.9	28.1	29.2	29.9	28.1	
COD	<20	<20	<20	<20	<20		
Fluoride	0.13	0.17	0.17	0.18			
Nitrate	6.24	3.54	5.49	6.42	5.91	<1	< 0.02
pH Upper Limit	8.2	7.7	7.2	7.5	7.3	8.1	7.3
pH Lower Limit							
Sulfate	178	178	179	171	165	169	
TOC	3.16	2.73	0.57	0.61			
TDS	440	435	435	422	420	422	
Benzene	<0.0050	<0.0050	<0.0050	<0.0050	<0.0050		
1,1-dichloroethape	< 0.0050	<0.0050	< 0.0050	<0.0030	< 0.0050		
1,1-dichloroethene	< 0.0050	<0.0050	<0.0050	< 0.0050	< 0.0050		
Tolucne	< 0.0050	< 0.0050	< 0.0050	< 0.0050	<0.0050		
1,1,1-trichloroethane	< 0.0050	< 0.0050	<0.0050	< 0.0050	< 0.0050		
	All concentr	ations in mg/l	4.				

Dixon Marquette Cement Comp Background MW1-S Statistical Analysis

1,1-dichloroethane0.005Method Detection Limit1,1-dichloroethene0.005Method Detection LimitToluene0.005Method Detection Limit	atisticai Analysis	Sample 14	Mean	Stnd. dev.	Deg of Freedom	99 Percentile	Upper limit	Method
Arsenic   Barlum	Aluminum						0.040	Lowest non-detect
Barlum	Antimony							Test of proportions
Beryllium   Boron   Cadmium   Calcium   102   7   10   2.76   109   Student's t   Chromium   0.001   0.001   5   3.36   0.002   Cohen   Copper   0.003   0.002   5   3.36   0.003   Cohen   Copper   0.003   0.002   5   3.36   0.005   Cohen   Copper   0.003   0.002   5   3.36   0.005   Cohen   Copper   0.003   0.002   5   3.36   0.005   Cohen   Copper   0.003   0.005   2   3.36   0.005   Cohen   Copper   0.003   0.005   2   3.36   0.005   Cohen   Copper   0.003   0.005   2   3.36   0.005   Cohen   Copper   0.001   Lowest non-detect   Iron   0.236   0.075   2   3.36   0.065   Cohen   Copper   0.003   0.015   11   2.72   0.045   Student's t   Copper   0.003   0.003   5   3.36   0.007   Cohen   Co	Arsenic							Test of proportions
Boron   Cadmium   102   7   10   2.76   109   Student's t   Chrombum   0.001   0.001   5   3.36   0.002   Cohen   Cobalt   0.002   0.001   5   3.36   0.002   Cohen   Cobalt   0.002   0.001   5   3.36   0.003   Cohen   Copper   0.003   0.002   5   3.36   0.005   Cohen   Copper   0.003   0.002   5   3.36   0.005   Cohen   Copper   0.003   0.002   5   3.36   0.005   Cohen   Comper   0.003   0.005   Cohen   Cohen   Cobalt   Cohen   Cohe	Barium		0.073	0.021	10	2.76	0.091	Student's t
Cadmium	Beryllium			•			0.001	Lowest non-detect
Calcium	Boron							Test of proportions
Chromium	Cadmium						0.001	Lowest non-detect
Cobalt	Calcium		102	7	10	2.76	109	Student's t
Cohalt	Chromium		0.001	0.001	5		0.002	Cohen
Lead	Cobalt		0.002	0.001	5	3.36 *-		Cohen
Iron	Copper		0.003	0.002	5	3.36	0.005	Cohen
Magnesium	Lead						0.001	Lowest non-detect
Manganese         0.033         0.015         11         2.72         0.045         Student's t           Mercury         0.0002         0.003         0.003         5         3.36         0.007         Cohen           Potassium         0.002         0.002         5         3.36         0.004         Cohen           Selenium         0.002         0.002         5         3.36         0.004         Cohen           Sodium         11.8         1.2         10         2.76         12.8         Student's i           Thallium         Test of proportions           Vanadium         Test of proportions           Zine         0.036         0.023         9         2.82         0.058         Student's t           Alkalinity, total         235         20         9         2.82         253         Student's t           Alkalinity, blearbonate         227         25         9         2.82         250         Student's t           Ammonia nitrogen         Test of proportions           Chloride         29.6         1.5         11         2.72         30.8         Student's t           COD         5.0         Lowest non-detec	Iron		0.236	0.075	2	3.36	0.465	Cohen
Mercury	Magnesium		50.2	3.9	9	2.82	53.9	Student's t
Nickel   0,003   0,003   5   3,36   0,007   Cohen   Test of proportions	Manganese		0.033	0.015	11	2.72	0.045	Student's t
Potassium   Selentum   0.002   0.002   5   3.36   0.004   Cohen   Sodium   11.8   1.2   10   2.76   12.8   Student's i   Test of proportions   Test of p	Mercury						0.0002	Lowest non-detect
Scientum	Nickel		0.003	0.003	5	3.36	0.007	Cohen
Sodium	Potassium							Test of proportions
Thallium   Test of proportions   Test of proportions   Test of proportions   Test of proportions   Test of proportions   Test of proportions   Test of proportions   Test of proportions   Student's t	Selenium		0.002	0.002	5	3.36	0.004	Cohen
Vanadium   Test of proportions   Zinc   O.036   O.023   9   2.82   O.058   Student's t	Sodium		11.8	1.2	10	2.76	12.8	Student's t
Alkalinity, total   235   20   9   2.82   253   Student's t	Thallium							Test of proportions
Alkalinity, total       235       20       9       2.82 section       253       Student's tend for proportions         Alkalinity, phenol.       - East of proportions         Alkalinity, blearbonate       227       25       9       2.82       250       Student's tend for proportions         Chloride       29.6       1.5       11       2.72       30.8       Student's tend for proportions         COD       5.0       Lowest non-detect         Fluoride       0.08       0.09       5       3.36       0.19       Cohen         Nitrate       7.32       1.26       5       3.36       11.07       Cohen         pil Upper Limit       7.516       0.338       12       3.06       7.82       Student's t         pil Lower Limit       7.52       0.338       12       3.06       7.22       Student's t         Sulfate       170.8       9.2       11       2.72       178.3       Student's t         TOC       2.7       1.9       9       2.82       4.5       Student's t         TDS       484       69       11       2.72       540       Student's t         Benzene       0.005       Method Detection Limit       <	Vanadium							Test of proportions
Alkalinity, phenol.         Test of proportions           Alkalinity, bicarbonate         227         25         9         2.82         250         Student's t           Ammonia nitrogen         Test of proportions           Chloride         29.6         1.5         11         2.72         30.8         Student's t           COD         5.0         Lowest non-detect         5.0         Lowest non-detect           Fluoride         0.08         0.09         5         3.36         0.19         Cohen           Nitrate         7.32         1.26         5         3.36         11.07         Cohen           pil Upper Limit         7.516         0.338         12         3.06         7.82         Student's t           pll Lower Limit         7.52         0.338         12         3.06         7.22         Student's t           Sulfate         170.8         9.2         11         2.72         178.3         Student's t           TOC         2.7         1.9         9         2.82         4.5         Student's t           TDS         484         69         11         2.72         540         Student's t           Benzene         0.005 </td <td>Zinc</td> <td></td> <td>0.036</td> <td>0.023</td> <td>9</td> <td>2.82</td> <td>0.058</td> <td>Student's t</td>	Zinc		0.036	0.023	9	2.82	0.058	Student's t
Alkalinity, bicarbonate         227         25         9         2.82         250         Student's t           Ammonia nitrogen         Test of propertions           Chloride         29.6         1.5         11         2.72         30.8         Student's t           COD         5.0         Lowest non-detect           Fluoride         0.08         0.09         5         3.36         0.19         Cohen           Nitrate         7.32         1.26         5         3.36         11.07         Cohen           pH Upper Limit         7.516         0.338         12         3.06         7.82         Student's t           pH Lower Limit         7.52         0.338         12         3.06         7.22         Student's t           Sulfate         170.8         9.2         11         2.72         178.3         Student's t           TOC         2.7         1.9         9         2.82         4.5         Student's t           TDS         484         69         11         2.72         540         Student's t           Benzene         0.005         Method Detection Limit         0.005         Method Detection Limit           1,1-dichloroethene	Alkalinity, total		235	20	9	2.82	253	Student's t
Test of propertions	Alkalinity, phenol.					•		Test of proportions
Chloride         29.6         1.5         11         2.72         30.8         Student's t           COD         5.0         Lowest non-detect           Fluoride         0.08         0.09         5         3.36         0.19         Cohen           Nitrate         7.32         1.26         5         3.36         11.07         Cohen           pH Upper Limit         7.516         0.338         12         3.06         7.82         Student's t           pH Lower Limit         7.52         0.338         12         3.06         7.22         Student's t           Sulfate         170.8         9.2         11         2.72         178.3         Student's t           TOC         2.7         1.9         9         2.82         4.5         Student's t           TDS         484         69         11         2.72         540         Student's t           Benzene           1,1-dichloroethane         0.005         Method Detection Limit           1,1-dichloroethene         0.005         Method Detection Limit           Toluene         0.005         Method Detection Limit	Alkalinity, bicarbonate		227	25	9	2.82	250	Student's t
COD   S.0   Lowest non-detect	Ammonia nitrogen							Test of propertions
Fluoride	Chloride		29.6	1.5	11	2.72	30.8	Student's t
Nitrate   7.32   1.26   5   3.36   11.07   Cohen	COD						5.0	Lowest non-detect
pH Upper Limit         7.516         0.338         12         3.06         7.82         Student's t           pH Lower Limit         7.52         0.338         12         3.06         7.22         Student's t           Sulfate         170.8         9.2         11         2.72         178.3         Student's t           TOC         2.7         1.9         9         2.82         4.5         Student's t           TDS         484         69         11         2.72         540         Student's t           Benzene           1,1-dichloroethane         0.005         Method Detection Limit           1,1-dichloroethene         0.005         Method Detection Limit           Toluene         0.005         Method Detection Limit	Fluoride		0.08	0.09		3.36	0.19	Cohen
pH Lower Limit         7.52         0.338         12         3.06         7.22         Student's t           Sulfate         170.8         9.2         11         2.72         178.3         Student's t           TOC         2.7         1.9         9         2.82         4.5         Student's t           TDS         484         69         11         2.72         540         Student's t           Benzene           1,1-dichloroethane         0.005         Method Detection Limit           1,1-dichloroethene         0.005         Method Detection Limit           Toluene         0.005         Method Detection Limit	Nitrate		7.32	1.26	5	3.36	11.07	Cohen
Sulfate	pH Upper Limit		7.516	0.338	12	3.06	7.82	Student's t
TOC   2.7   1.9   9   2.82   4.5   Student's t	pH Lower Li <b>mit</b>		7.52	0.338	12	3.06	7.22	Student's t
### TDS   484   69   11   2.72   540   Student's t      Benzene	Sulfate		170.8	9.2		2.72	178.3	Student's t
Benzene 0.005 Method Detection Limit 1,1-dichloroethane 0.005 Method Detection Limit 1,1-dichloroethene 0.005 Method Detection Limit Toluene 0.005 Method Detection Limit O.005 Method Detection Limit	TOC		2.7	1.9	9	2.82	4.5	Student's t
1,1-dichloroethane 0.005 Method Detection Limit 1,1-dichloroethene 0.005 Method Detection Limit Toluene 0.005 Method Detection Limit O.005 Method Detection Limit	TDS		484	69	11	2.72	540	Student's t
1,1-dichloroethene 0.005 Method Detection Limit Toluene 0.005 Method Detection Limit	Benzene							Method Detection Limit
Tolucne 0.005 Method Detection Limit	1,1-dichloroethane						0.005	Method Detection Limit
	1,1-dichloroethene						0.005	Method Detection Limit
1,1,1-trichlorocthane 0.005 Method Detection Limit	Toluene							Method Detection Limit
	1,1,1-trichloroethane						0.005	Method Detection Limit

All concentrations in mg/L.

Monitoring Well 1-WT	Monitoring Well 5-S	Zq	Monitoring Well 6-S Zd	Veli 6-S	pZ	Monitoring Well 7-S		Zd Mon	Monitoring Well 8-5	ΡZ
Detected Sample Size	Samples Detected Sample Size	57	Samples Detected	Sample Size	<del>-1</del>	Samples Detected Samp	le Size	Samples	Detected Sample Si	çe Ze
9	2 2	0.91	0	7	0.91		7 0	) 61.	7	16.0
01 10	7	0.99	'n	7	2.61		7 3	ý (99·	-	2,10
3 10	6 7	7.27	9		2.27	9	7 2			2.89
6 12	7 7	2.26	7	7	2.26	7	7 2	.26		2.26
6	0 2	16.0	•	7	0.19	0	7 0	.91	7	1.45
9	6 7	1.69	• 9	-	1.69	7	7 2	.38	7	1.08
9	0	0.86	0	4	98.0	6	4	.84	4	98.0
2 10	7	0.36		7	0.30		7 0	30	7	1.26
	0 7	0.82	0	1	0.82	0	7 0	.82	7	0.82
	Detected Sample Size 9 10 10 5 12 6 12 9 11	Detected Sample Size  9	c Samples Detected Sample Size  2 7 0.99  6 7 2.26  7 2.26  0 7 0.99  6 7 0.99  6 7 0.91  6 7 0.91  7 0.91  6 7 0.91  7 0.92  0 7 0.91  0 7 0.93	Detected Sample Size         Samples Detected Sample Size         Samples Detected           9         0         7         0.91         0           10         2         7         0.99         5           12         7         2.26         7         6           9         6         7         0.91         1         6           9         6         7         1.69         6         6         6         6           10         1         7         0.36         1         6         1         1         6         1         1         6         1         1         1         0         7         0.36         1         1         0         1         0         0         2         0	Detected Sample Size         Sample Size         Sample Size         Sample Size           9         0         7         0.91         0         7           10         2         7         0.99         5         7           10         6         7         2.27         6         7           12         7         7         2.26         7         7           9         6         7         1.69         6         7           1         9         6         7         1.69         6         7           1         10         1         7         0.36         1         7           11         0         7         0.36         1         7         7	Detected Sample Size         Samples Detected Sample Size         Samples Detected Sample Size         1           9         0         7         0.91         0         7         0.91           10         6         7         2.26         7         2.27           5         12         7         7         2.26         7         2.27           9         6         7         0.91         1         0         7         0.19           9         6         7         1.69         6         7         1.69           6         0         4         0.36         1         7         0.30           10         1         7         0.36         1         7         0.30           11         0         7         0.82         0         7         0.82	Detected Sample Size         Samples Detected Sample Size         Samples Detected Sample Size         Samples Detected Sample Size         Samples Detected Samples Det	Samples Detected Sample Size         Samples Detected Sample Size           0         7         0.91         1         \$ 7         7         6         7	Samples Detected Sample Size         Samples Detected Sample Size           0         7         0.91         1         4         7         0.19           5         7         2.61         7         3.66         7         2.27           6         7         2.26         7         7         2.26           1         9         7         0.19         0         7         0.91           6         7         1.69         7         7         2.38           0         4         0.86         3         4         1.84           1         7         0.30         1         0         0         0           0         7         0.82         0         7         0.30	Samples Detected Sample Size         Samples Detected Samples Size         Samples Detected Sample Size         Samples Detected Samples Size         Samples Detected Sample

If Zd > 1.96, then sample set is significantly different than background

	Sample 1	Sample 2	Sample 3	Sample 4	Sample 5	Sample 6	Sample 7
Aluminum	0.27	0.08	0.226	0.074	< 0.04	< 0.04	< 0.04
Antimony	< 0.001	< 0.001	< 0.001	< 0.001	< 0.001	< 0.001	< 0.001
Arsenic	0.003	0.002	0.001	0.002	0.002	0.001	0.0005
Barium	0.122	0.155	0.14	0.183	0.25	0.202	0.142
Beryllium	< 0.001	< 0.001	< 0.001	< 0.001	< 0.001	< 0.001	< 0.001
Boron	< 0.05	< 0.05	0.061	0.078	< 0.05	< 0.05	0.031
Cadmium	< 0.001	< 0.001	< 0.001	< 0.001	< 0.001	< 0.001	< 0.001
Calcium	97. <b>7</b>	107	99.1	110	96.4	129	125
Chromium	0.002	0.001	0.001	0.001	0.001	< 0.001	< 0.001
Cobalt	0.007	0.008	0.001	0.004	0.004	0.005	0.005
Copper	0.012	0.011	0.005	0.005	0.001	0.008	0.004
Lead	0.003	0.004	0.001	0.001	< 0.001	< 0.001	< 0.001
Iron	0.73	0.501	0.175	0.966	0.304	0.62	0.582
Magnesium	18.2	32.7	35.7	39.6	38.8	49.4	46.8
Manganese	2.04	2.64	0.912	2.23	1.88	1.78	2.36
Mercury	< 0.0002	< 0.0002	<0.0002	<0 0002	< 0.0002	0.0002	< 0.0002
Nickel	0.023	0.016	0.009	0.01	0.022	0.018	0.017
Potassium	2.43	3.06	3.63	1.9	3.3	4	3.3
Selenium	< 0.004	< 0.001	<0.001	< 0.001	. <0.001	0.005	0.002
Sodiu <b>m</b>	64.9	38.5	35.1	34.4	26.2	31.5	26.1
Thallium	< 0.001	< 0.001	< 0.001	0.002	< 0.001	0.001	< 0.001
Vanadium	0.005	0.002	0.002	0.007	0.004	0.007	0.0005
Zinc	0.01	0.011	0.008	0.012	0.001	0.014	0.006
Alkalinity, total	201	346	322	215	270	290	280
Alkalinity, phenol	<1.0	<1.0	<1.0				
Alkalinity, bicarbonate	198	330	320	188	265	276	272
Ammonia nitrogen	<0.1	<0.1	<0.1	<0.1	<0.1	0.14	<0.01
Chloride	1.55	22.1	17.9	18.8	18.2	19.6	18.6
COD	<20	44	<20	25.6	<20	42.9	<20
Fluoride	0.13	1.13	0.12	0.1	0.13	0.13	0.15
Nitrate	<0.1	<0.1	0.12	<0.1	<0.1	<0.1	<0.1
pH Upper Limit	7.5	8.3	7.8	8	7.4	7.1	7.6
pH Lower Limit	04.0	0.6.1					
Sulfate	83.8	86.1	95.7	120	175	238	253
TOC	8.06	22.3	4.04	7.44	10.4	1.28	1.43
TDS	320	430	407	390	420	485	471
Benzene	<0.0050	<0.0050	< 0.0050	<0.0050	< 0.0050	<0.0050	<0.0050
1,1-dichloroethane	< 0.0050	< 0.0050	<0.0050	< 0.0050	<0.0050	< 0.0050	< 0.0050
1,1-dichloroethene	< 0.0050	< 0.0050	< 0.0050	< 0.0050	<0.0050	< 0.0050	< 0.0050
Toluene	< 0.0050	< 0.0050	< 0.0050	< 0.0050	· <0.0050	<0.0050	<0.0050
1,1,1-trichloroethane	< 0.0050	< 0.0050	< 0.0050	< 0.0050	<0.0050	< 0.0050	< 0.0050
	All concentr	ations in mg	/L.		-		

•	Mean	Stnd, dev.	Deg of Freedom	99 Percentile	Upper limit	Method
Aluminum	0.163	0.100	3	4.54	0.454	Cohen
Antimony					0.001	Lowest non-detect
Arsenic	0.002	0.001	6	3.14	0.003	Student's t
Barium	C.171	0.044	6	3.14	0.227	Student's t
Beryllium					0.001	Lowest non-detect
Boron						Test of proportions
Cadmium					0.0	Lowest non-detect
Calcium	109.2	13.2	6	3.14	126.1	Student's t
Chromium	0.001	0.000	· 4	3.75	0.002	Cohen
Cobalt	0.005	0.002	6	3.14	0.008	Student's t
Copper	0.007	0.004	6	3.14	0.012	Student's t
Lead	0.002	0.002	3	4.54	0.006	Cohen
Iron	0.6	0.3	6	3.14	0.9	Student's t
Magnesium	37.3	10.3	6	3.14	50.5	Student's t
Manganese	1.98	0.55	6	3.14	2.69	Student's t
Mercury						Test of proportions
Nickel	0.0	0.0	6	3.14	0.0	Student's t
Potassium	3.1	0.7	6	3.14	4.0	Student's t
Selenium						Test of proportions
Sodium	36.7	13.3	6	3.14	53.7	Student's t
Thallium						Test of proportions
Vanadium	0.004	0.603	6	3.14	0.007	Student's t
Zinc	0.009	0.004	6	3.14	0.014	Student's t
Alkalinity, total	275	53	6	3.14	342	Student's t
Alkalinity, phenol					1	Lowest non-detect
Alkalinity, bicarbonate	264	55	6	3.14	334	Student's t
Ammonia nitrogen						Test of proportions
Chloride	16.7	6.8	6	3.14	25.4	Student's t
COD						Test of proportions
Fluoride	0.27	0.38	6	3.14	0.76	Student's t
Nitrate			•			Test of proportions
pH Upper Limit	7.65	0.37	7	3.5	8.14	Student's t
pH Lower Limit	7.65	0.37	7	3.5	7.16	Student's t
Sulfate	150.2	72.2	6	3.14	242.8	Student's t
TOC	8	7	6	3.14	17	Student's t
TDS	418	55	6	3.14	488	Student's t
Benzene					0.005	Method Detection Limit
1,1-dichloroethane					0.005	Method Detection Limit
1,1-dichloroethene					0.005	Method Detection Limit
Toluene					0.005	Method Detection Limit
1,1,1-trichloroethane					0.005	Method Detection Limit
	All conce	entrations in m	g/L.			

Background Tables

Dixon Marquette Cement Comp	TW-1WR	Test of Proportions
Dix	7	Ž

	Me stocing Well I-WT	CLI I-WT	Mositoriae Well S-S		P2	Monitoring Well 6-S	cli 6-S	P7	Monitoring Well 7-S	117.S	PZ	Monitoring Well 8-S	S-8-13
	Samuelos Detected Sample Size Samples	Sample Size	Samples Detected 5	.5	-	Samples Detected Sample Size	ample Size		Samples Detected Sample Size	ample Size		Samples Detected Sample Size	ample Si
Boron	~	,	9	7		9	-	1.67	9	-	16.	7	_
Mercury		r~	0	1~	8.	و.	~	3	0	7	3.	0	7
Selenium	- (1	-	··c	۲	2,16	ý	۲	2.16	9	7	2.16	7	7
Thailium	. 61	-1		~	1.53		~	0.65	0	7	1.53	m	7
Ammonia nitrosen		r~	ı emi	^	0.00		۲-	0.0	ent	7	0.00	0	2
, aoo	. ~	(	• •	~	0.53	m	1~	0.0	14	7	0.56	71	7
N. C.		۲۰	~	t-	8	7	7	8	c	7	40.	1	7

If Zd > 1.96, then sample set is significantly different than background

	Sample 1	Sample 2	Sample 3	Sample 4	Sample 5	Sample 6
Aluminum	0.08	0.109	<0.1	<0.1	< 0.04	0.111
Antimony	< 0.1	< 0.1	< 0.001	< 0.001	< 0.001	< 0.001
Arsenic	< 0.005	< 0.08	<0.08	<0.08	< 0.001	< 0.001
Barium	0.063	0.044	0.047	0.046	0.047	0.053
Beryllium	<0.01	<0.01	< 0.01	< 0.001	< 0.001	< 0.001
Boron	<0.1	< 0.05	< 0.05	<0.05	<0.05	< 0.05
Cadmium	< 0.005	< 0.02	< 0.02	< 0.02	<0.02	< 0.001
Calcium	93	80	86	79	92.3	87.6
Chromium	0.002	0.002	< 0.04	< 0.02	< 0.02	< 0.02
Cobalt	0.004	0.003	< 0.02	< 0.02	< 0.02	<0.001
Copper	0.002	0.003	0.004	0.003	0.007	0.001
Lead	< 0.005	<0.1	< 0.1	<0.01	< 0.001	< 0.001
Iron	0.872	0.79	0.273	0.568	0.519	1.19
Magnesium	39	44	41	49.7	45.5	44
Manganese	0.071	0.066	0.007	0.058	0.053	0.001
Mercury	< 0.0002	< 0.0002	< 0.0002	< 0.0002	< 0.0002	< 0.0002
Nickel	0.002	0.003	0.003	0.003	0.002	0.001
Potassium	1.5	1.4	1.57	1.34	1.89	2.9
Selenium	100.0	. 0.001	0.004	0.001	0.006	0.004
Sedlum	3.1	3.8	3.1	4.2	3.3	3.65
Thallium	<1	<1	< 0.001	< 0.001	< 0.001	0.002
Vanadium	0.002	0.002	0.001	0.002	0.004	0.001
Zinc	0.011	0.013	0.018	0.017	0.02	0.004
Alkalinity, total	348	350	350	282	268	320
Alkalinity, phenol.	<5	<5	<1	<1	<1	<1
Alkalinity, bicarbonate	348	350	350	250	250	310
Ammonia	<0.2	<0.2	<0.1	<0.1	0.11	< 0.1
Chloride	2.23	1.91	2.55	2.58	4.86	2.61
COD	<5	<5	<8.3	<20	<20	<20
Fluoride	0.18	0.13	0.14	0.11	0.14	0.13
Nitrate	0.2	0.18	0.19	0.14	0.13	0.2
pH Upper Limit	7.31	<b>7</b> .1	7.3	7.4	7.5	8.1
pH Lower Limit						
Sulfate	37	42	56	35	37	37.7
TOC	3.92	1.5	6	4.13	2.79	4.59
TDS	396	400	400	400	370	333
Benzene	<0.0005			< 0.0050	< 0.0050	<0.0050
1,1-dichloroethane	<0.0010	< 0.0050	<0.0010	<0.0050	< 0.0050	<0.0050
1,1-dichloroethene	< 0.0020	<0.0020	<0.0020	<0.0050	<b>&lt;0.0</b> 050	< 0.0050
Toluene	< 0.0010	<0.0010	<0.0010	<0.0050	< 0.0050	< 0.0050
1,1,1-trichloroethane	<0.0010	< 0.0010	<0.0010	<0.0050	< 0.0050	< 0.0050
	All concent	rations in mg	y/L.,			

	Sample 7	Sample 8	Sample 9	Sample 10	Sample 11	Sample 12	Sample 13	Sample 14
Aluminum	<0.04	0.048	<0.04				•	
Antimony	< 0.001	< 0.001	< 0.001			•		
Arsenic	< 0.001	0.001	<0.001	. 0.001	< 0.001			
Barium	0.047	0.055	0.03	0.046	0.036			
Beryllium	< 0.001	< 0.001	< 0.001	< 0.001				
Boron	0.051	< 0.05	< 0.05	0.009				
Cadmium	< 0.001	< 0.001	< 0.001	< 0.001	< 0.001	< 0.001		
Calcium	84.9	89.4	51.8	82.4	69.1			
Chromium	< 0.001	0.001	<0.001	0.062	< 0.001			
Cobalt	0.007	< 0.001	< 0.001	10 ).0>	+		•	
Copper	<0.02	< 0.02	< 0.02	<0.001				
Lead	< 0.001	0.001	0.002	< 0.001	< 0.001			
Iron	0.132	0.214	0.217	< 0.1	<0.1	<0.1		
Magnesium	121	36.2	43.7	38.5	٠.,	-0.1		
Manganese	0.219	0.002	0.001	< 0.01	< 0.01	< 0.01	< 0.01	<0.01
Mercury	< 0.0002	<0.0002	< 0.0002	< 0.0002	0.01	-0.01	V.01	10.0
Nickel	< 0.05	< 0.05	< 0.05	< 0.001				
Potassium	1.2	1.9	<2	<2	<2	<2	< 0.05	
Selentum	< 0.005	< 0.15	< 0.15	<0.15	<0.004	•	·().05	
Sodium	3.57	3.9	3.8	3	3.6	2.4		
Thallium	< 0.001	0.003	< 0.001	•	2.0	47.7		
Vanadium	< 0.05	< 0.05	< 0.02	< 0.001				
Zinc	<0.02	<0.02	<1	<0.001			•	
Alkalinity, total	314	325	312	328				
Alkalinity, phenol.							•	
Alkalinity, bicarbonate	296	308	310	315				
Ammonia	<0.1	0.29	<0.1					
Chloride	2.46	<5	<5	<5	<5		•	
COD	<20	52.6	45.7	<20		·		
Fluoride	0.17	<0.2	< 0.2	< 0.21				
Nitrate	0.18	<1	<1	<1	<1			
pH Upper Limit	8.1	7.7	7.2	7.4	7.2	8		
pli Lower Limit								
Sulfate	37	37.6	35.6	40.7	40.2	39.9		
TOC	3.16	0.28	<1	<0.5				
TDS	338	326	322	323	317	316		
Benzene	< 0.0050	<0.0050	<0.0050	< 0.0050				
1,1-dichloroethane	< 0.0050	< 0.0050	< 0.0050	< 0.0050				
1,1-dichloroethene	< 0.0050	< 0.0050	< 0.0050	< 0.0050				
Toluene	< 0.0050	< 0.0050	< 0.0050	< 0.0050				
1,1,1-trichloroethane	< 0.0050	<0.0050	< 0.0050	< 0.0050				
	All concentra	tions in mg/l	L.,					

Anaiysis	Mean	Stnd. dev.	Deg of Freedom	≠ 99 Percentile	Upper limit	
Aluminum						Test of proportions
Antimony					0.001	Lowest non-detect
Arsenic						Test of proportions
Barium	0.043	0.009	10	2.76	0.051	Student's t
Beryllium					0.001	Lowest non-detect
Boron	•	•				Test of proportions
Cadmium					0.001	Lowest non-detect
Calcium	81.4	11.9	10	2.76	91.8	Student's t
Chromium						Test of proportions
Cobalt						Test of proportions
Copper	0.003	0.002	5	3.36	<b>0.0</b> 06	Cohen
Lead						Test of proportions
Iron	0.531	0.361	8	2.9	0.819	Cohen
Magnesium	50.3	25.2	9	2.82	73.9	Student's t
Manganese	0.053	0.069	8	2.9	0.098	Cohen
Mercury					0.0002	Lowest non-detect
Nickel	0.002	0.001	5	3.36	0.003	Cohen
Potassium	1.7	0.5	7	3	2.3	Cohen
Selenium	0.003	0.002	5	3.36	0.005	Cohen
Sodium	3.45	0.49	11	2.72	3.85	Student's t
Thallium						Test of proportions
Vanadium	0.002	0.001	5	3.36	0.003	Cohen
Zinc	0.014	0.006	5	3.36	0.023	Cohen
Alkalinity, total	320	28	9	2.82	346	Student's t
Aikalinity, phenol.	320		-		1.000	Lowest non-detect
Alkalinity, bicarbonate	310	35	9	2.82	343	Student's t
Ammonia	2.0					Test of proportions
Chloride	2.74	0.97	6	3.14	4.1	Cohen
COD	_ · ·				5.0	Test of proportions
Fluoride	0.14	0.02	6	3.14	0.19	Cohen
Nitrate	0.17	0.03	6	3.14	0.21	Cohen
pH Upper Limit	7.526	0.362	11	3.11	7.86	Student's t
pli Lower Limit	7.53	0.362	11	3.11	7.19	Student's t
Sulfate	39.6	5.6	11	2.72	44.2	Student's t
TOC	3.3	1.8	7	3	5.1	Cohen
TDS	353	36	11	2.72	383	Student's t
Benzene					0.005	Method Detection Limit
1,1-dichloroethane					0.005	Method Detection Limit
1.1-dichloroethene					0.005	Method Detection Limit
Toluene					0.005	Method Detection Limit
1,1,1-trichloroethane					0.005	Method Detection Limit
1,1,1-ificulor octuanc			/1			

All concentrations in mg/L.

Adjusted Groundwater Standards Dixon Marquette Cement Company – Dixon, Illinois Prepared by Preston Engineering, Inc. February 27, 2001

DOWNGRADIENT TABLES

### Dixon Marquette Cement Company Monitoring Well 3S

Parameter	MW3-S	MW3-S	MW3-S	MW3-S	MW3-S	MW3-S
	07-May-98	17-Jun-98	24-Jul-98	26-Oct-98	16-Feb-99	26-Jul-99
Aluminum						0.29
Antimony					•	0.001
Arsenic	0.0176	-		0.0304		0.131
Barium	0.064		0.049	0.036		0.054
Beryllium						0.001
Boron						0.05
Cadmium	0.005			0.01		0.001
Calcium	8.9		4.4	3		4.44
Chromium	0.04			0.04		0.003
Cobalt						0.001
Copper						0.011
Lead	0.005					0.001
Iron	1.46	1.0	1.0	0.1		0.484
Magnesium						4.07
Manganese	0.051	0.1	0.073	0.031		0.101
Mercury					•	0.0002
Nickel						0.014
Potassium	1400	1600	1600	1900	200	2314
Selenium	0.005					<b>0</b> .006
Sodium	82	100	98	120	120	145
Thallium						0.001
Vanadium						0.015
Zinc						0.008
Alkalinity, total	980		1400	1500		1590
Alkalinity, phenol.			640	650		
Alkalinity, bicarb	240		120	200		1450
Ammonia nitrogen						0.10
Chloride	137	130	160	.140	170	181
COD						20
Fluoride						3.97
Nitrate as N	1.0	1.0	1.0	1.0		0.10
pli Upper Limit (stnd units)	10.2	10.4	10.5	10.4	10.4	10.9
pli Lower Limit (stnd units)						
Sulfate	690	1200	1400	980	1100	980
TOC						12.9
Ttl Dis Solids	2960	3700	3900	4200		3776
Benzene	0.001		0.0005	0.0005		0.005
1,1-dichloroethane	0.0029	0.003		0.004		0.005
1,1-dichloroethene	0.001		0.002			0.005
Toluene	0.001		0.001	0.001		0. <b>005</b>
1,1,1-trichloroethane	0.015	0.0126	0.0122	0.0114		0.005

All values in mg/L unless noted otherwise

XXX Non detected Concentration

XXX Concentration > Average + 1 Standard Deviation

Parameter	MW3-S	MW3-S	MW3-S	MW3-S	MW3-S	99.5% CI*
	<b>09-</b> Sep-99	01-Nov-99	02-Mar-00	23-Jun-00	29-Sep-00	29-Nov-00
Aluminum	0.39	0.837	0.406	0.071	0.233	0.028
Antimony	0.001	0.001	0.001	0.001	0.001	0.001
Arsenic	0.079	0.091	0.065	0.06	0.097	0.175
Barium	0.046	0.051	0.042	0.034	0.042	0.089
Beryllium	0.001	0.001	0.001	0.001	0.001	0.001
Boron	0.15	0.062	0.184	0.05	0.052	0.204
Cadmium	0.001	0.001	0.001	0.001	0.001	0.001
Calcium	3.71	5.34	2.4	3.53	6.1	1.3
Chromium	0.001	0.003	0.001	0.003	0.003	0.018
Cobalt	0.002	0.003	0.001	0.001	0.001	0.001
Copper	0.013	0.014	0.009	0.006	0.045	0.022
Lead	0.002	0.002	0.001	0.001	0.001	0.001
Iron	0.399	0.861	0.18	0.152	0.291	0.409
Magnesium	2.8	3.43	1.1	2.91	3.5	0.6
Manganese	0.173	0.328	0.106	0.001	0.056	0.041
Mercury	0.0002	0.0002	0.0002	0.0002	0.0002	0.0015
Nickel	0.013	0.015	0.008	0.008	0.013	0.038
Potassium	2010	2170	2120	1980	218	2120
Selenium	0.005	0.007	0.004	0.004	0.005	0.016
Sodium	139	134	130	130	138	129
Thallium	0.001	0.001	0.001	0.001	0.002	0.001
Vanadium	0.001	0.011	0.003	0.01	0.016	0.029
Zinc	0.012	0.014	0.012	0.001	0.014	0.015
Alkalinity, total	1600	1472	1540	1507	1630	1660
Alkalinity, phenol.			662	1155	710	774
Alkalinity, bicarb	250	800	293	328	807	746
Ammonia nitrogen	1.60	0.10	0.10	0.10	0.61	0.10
Chloride	163	151	160	231	202	203
COD	35	20	36.4	242	68.6	55.9
Fluoride	4	4.44	3.47	3.11	3.02	3.2
Nitrate as N	0.10	0.16	0.10	0.10	0.10	0.10
pH Upper Limit (stnd units)	11.2	10.5	10.8	10.8	10.4	10.7
pH Lower Limit (stnd units)						
Sulfate	1054	1009	931	1153	1170	1242
тос	22.7	17	27.2	17.9	9.26	14.3
Ttl Dis Solids	3462	4350	3980	3150	4400	4390
Benzene	0.005	0.005	0.005	0.005	0.005	0.005
1,1-dichloroethane	0.005	0.005	0.005	0.005	0.005	0.005
1,1-dichloroethene	0.005	0.005	0.005	0.005	0.005	0.005
Toluene	0.005	0.005	0.005	0.005	0.005	0.005
1,1,1-trichloroethane	0.0068	0.0064	0.005	0.005	0.005	0.005

All values in mg/L unless noted otherwise

XXX Non detected Concentration

XXX Concentration > Average + 1 Standard Deviation

## Dixon Marquette Cement Company Monitoring Well 3S

Parameter .	Average	Stnd Dev	Average + 1 Stnd Dev	Average + 2 Stnd Dev	99.5% CI*	Ave + 99.5% CI*
	30-Mar-01	30-Mar-01	30-Mar-01	30-Mar-01	30-Mar-01	30-Mar-01
Aluminum	0.322	0.269	0.591	0.861	0.408	0.730
Antimony	0.001	0.000	0.001	0.001	0.000	0.001
Arsenic	0.083	0.049	0.132	0.180	0.058	0.141
Barium	0.051	0.016	0.067	0.083	0.017	0.068
Beryllium	0.001	0.000	0.001	0.001	0.000	0.001
Boron	0.107	0.069	0.177	0.246	0.105	0.212
Cadmium	0.002	0.003	0.006	0.009	0.004	0.006
Calcium	. 4.3	2.1	6.4	8.6	2.3	6.6
Chromium	0.012	0.016	0.029	0.045	0.020	0.032
Cobalt	0.001	0.001	0.002	0.003	0.001	0.003
Copper	0.017	0,013	0.030	0.044	0.020	0.037
Lead	0.002	0.001	0.003	0.005	0.002	0.004
Iron	0.576	0.440	1.016	1.456	0.441	1.017
Magnesium	2.6	1.3	3.9	5.2	2.0	4.6
Manganese	0.096	0.090	0.186	0.276	0.090	0.186
Mercury	0.0004	0.0005	0.0009	0.0014	0.0007	0.0011
Nickel	0.016	0.010	0.026	0.036	0.016	0.031
Potassium	1636	718	2354	3072	673	2309
Selen'um	0.007	0.004	0.010	0.014	0.005	0.012
Sodium	122.1	19.2	141.3	160.5	18.4	140.5
Thallium	0.001	0.000	0.002	. 0.002	0.001	0.002
Vanadium	0.012	0.009	0.021	0.031	0.014	0.026
Zinc	• 0.011	0.005	0.016	0.021	0.007	0.018
Alkalinity, total	1488	195	1683	1878	211	1699
Alkalinity, phenol.	765.2	197.3	962.5	1159.8	355.6	1120.8
Alkalinity, bicarb	523	419	942	1361	454	977
Ammonia nitrogen	0.39	0.57	0.95	1.52	0.86	1.25
Chloride	169.0	30.4	199.4	229.7	28.5	197.5
COD	68.3	78.7	146.9	225.6	119.1	187.4
Fluoride	3.60	0.54	4.14	4.68	0.82	4.42
Nitrate as N	0.43	0.45	0.88	1.33	0.45	0.88
pH Upper Limit (stnd units)	10.6	0.3	10.9	11.2	0.3	:0.9
pH Lower Limit (stnd units)	4000					10.3
Sulfate	1076	180	1256	1436	169	1245
TOC	17.32	6.06	23.39	29.45	9.18	26.51
Ttl Dis Solids	3843	495	4337	4832	496	4338
Benzene	NA	NA	NA	NA	NA	NA
1,1-dichloroethane	0.0045	0.001	0.005	0.006	0.0009	0.0054
1,1-dichloroethcae	NA	NA	NA	NA	NA	NA
Toluene	NA	NA	NA	NA	NA	NA.
1,1,1-trichloroethane	0.0081	0.0038	0.012	· 0.016	0.0039	0.012

All values in mg/L unless noted otherwise

NA All sampling events below detection limit

<sup>\* 95%</sup> CI computed for all parameter except pH. A two sided 99% CI was computed.

### Dixon Marquette Cement Company Monitoring Well 5S

Parameter	MW5-S						
	26-Jul-99	09-Sep-99	01-Nov-99	02-Mar-00	23-Jun-00	29-Sep-00	29-Nov-00
Aluminum	0.17	0.17	0.161	0.286	0.041	0.151	0.195
Antimony	0.001	0.001	0.001	0.001	0.001	0.001	0.001
Arsenic	0.001	0.001	0.002	(元,0.005)一型	0.001	0.001	0.001
Barium	0.199	0.069	0.095	0.307	0.078	0.131	0.087
Beryllium	0.001	0.001	0.001	0.001	0.001	0.001	0.001
Boron	0.52	0.55	0.527	0.700	0.05	0.572	0.328
Cadmium	0.001	0.001	0.001	0.002	0.001	0.001	0.001
Calcium	274	291	279	278	204	177	175
Chromium	0.004	0.001	0.003	0.006	0.002	0.004	0.002
Cobalt	<b>0</b> .001	C.001	0.001	0.006	0.001	0.001	0.001
Copper	0.004	0.006	0.007	0.017	0.001	0.003	0.004
Lead	0.001	0.001	0.001	0.001	0.001	0.001	0.001
Iron	0.913	0.816	0.718	2.24	0.351	0.604	0.646
Magnesium	89	34.6	101	202	28.9	58.7	62.1
Manganese	0.081	0.064	0.047	1.25	0.001	0.044	0.117
Mercury	0.0002	0.0002	0.0002	0.0002	0.0002	0.0002	0.0002
Nickel	0.005	0.004	0.003	0.021	0.001	0.002	0.002
Potassium	23.5	61.9	59.5	341	23.9	74.5	17.7
Selenium	0.003	0.001	0.004	0.006	0.001	0.001	0.001
Sodium	22.8	78.5	55.3	34	47.2	18.6	9
Thallium	0.001	0.001	0.001	0.001	0.001	0.001	0.001
Vanadium	0.007	0.001	0.005	0.023	0.005	0.007	0.003
Zinc .	0.011	0.02	0.012	250.047	0.001	0.009	0.016
Alkalinity, total	324	372	586	386	432	490	502
Alkalinity, phenol.				1.0	1.0	1.0	1.0
Alkalinity, bicarb	314	360	540	377	427	375	479
Ammonia nitrogen Chloride	0.10	0.10	0.10	0.10	0.10	0.19	
COD	21.6	111	95.5	125	36.9	25.2	23.9
Fluoride	20	26	20	20	57.9		203
Nitrate as N	0.13 1.31	0.06	0.09 2.00	0.06	0.13	0.13	0.12
pH Upper Limit (stnd units)	7.6	7.8	7.6	0.10 6.9	0.10	0.10	0.10
pH Lower Limit (stnd units)	7.0	7.0	7.0	0.9	7.1	7.3	7.7
Sulfate	220	571	578	650	278	196	203
TOC	21.6	29	17.9	0.50	9.86	1.41	
Ttl Dis Solids	668	1090	1016	1213	654	633	1.46 646
Ren : ve	0.005	0.005					
Benane 1,1-dichloroethane	0.005 0.005	0.005	0.005	0.005	0.005	0.005	0.005
1,1-dichloroethene	0.005 0.005	0.005 0.005	0.005	0.005	0.005	0.005	0.005
Toluene	0.005 0.005	0.005 0.065	0.505 0.005	0.005 0.005	0.005	0.005	0.005
1,1,1-trichloroethane	0.005	0.005	0.005	0.005	0.005 0.005	0.005 0.005	0.00 <b>5</b> 0.00 <b>5</b>
							<del></del>

All values in mg/L unless noted otherwise

XXX Non detected Concentration

XXX Concentration > Average + 1 Standard Deviation

## Dixon Marquette Cement Company Monitoring Well 5S

Parameter	Average	Stnd Dev	Average + 1 Stnd Dev	Average + 2 Stnd Dev	99.5% CI*	Ave+ 99.5% CI*	Slope of Trend
	30-Mar-01	30-Mar-01	30-Mar-01	30-Mar-01	30-Mar-01	30-Mar-01	26-Jan-01
Aluminum	0.168	0.072	0.240	0.312	0.109	0.277	-0.00005
Antimony	0.001	0.000	0.001	0.001	0.000	0.001	
Arsenic	0.002	0.001	0.003	0.005	0.002	0.004	<b>-0</b> .0000 <b>008</b>
Barium	0.138	0.087	0.225	0.311	0.131	0.269	-0.00008
Beryllium	0.001	0.000	0.001	0.001	0.000	0.001	
Boron	0.464	0.213	0.677	0.889	0.322	0.786	-0.0004
Cadmium	0.001	0.000	0.002	0.002	0.001	0.002	-5E-08
Calcium	240	52	292	344	79	318	-0.25505
Chromium	0.003	0.002	0.005	0.006	0.003	0.006	-0.0000002
Cobalt	0.002	0.002	0.004	0.005	0.003	0.005	-0.0000003
Copper	0.006	0.005	0.011	0.016	0.008	0.014	-0.000006
Lead	0.001	0.000	0.001	0.001	0.000	0.001	
Iron	0.898	0.618	1.516	2.134	0.936	1.834	-0.0007
Magnesium	82.3	58.9	141.2	200.1	89.2	171.5	-0.0531
Manganese	0.229	0.452	0.681	1.132	0.684	0.913	-0.00005
Mercury	0.0002	0.0000	0.0002	0.0002	0.0000	0.0002	0.00003
Nickel	0.005	0.007	0.012	0.019	0.011	0.016	-0.000007
Potassium	86.0	114.7	200.7	315.3	173.6	259.6	-0.0304
Selenium	0.002	0.002	0.004	0.006	0.003	0.005	-0.000004
Sodium	37.9	24.1	62.0	86.1	36.5	74.4	-0.0749
Thallium	0.001	0.000	0.001	0.001	0.000	0.001	-0.0747
Vanadium	0.007	0.007	0.015	0.022	0.011	0.001	-0.0000002
Zinc	0.017	0.015	0.031	0.046	0.022	0.039	-0.000002
Alkalinity, total	442	90	532	622	136	578	0.205
Alkalinity, phenol.	1.0	0	1.0	1.0	0	1	
Alkalinity, bicarb	410	77	488	565	117	528	0.1217
Ammonia nitrogen	0.11	0.03	0.15	0.18	0.05	0.16	0.00008
Chloride	62.7	45.7	108.5	154.2	69.3	132.0	-0.1106
COD	53.2	67.4	120.6	188.0	102.1	155.3	0.23
Fluoride	0.10	6.03	0.14	0.17	0.05	0.15	0.001
Nitrate as N	1.42	2.26	3.68	5.93	3.42	4.84	-0.0075
pH Upper Limit (stnd units)	7.4	0.3	7.8	. 8.1	0.4	7.9	-0.0005
pH Lower Limit (stnd units)	206	204	500			7.0	
Sulfate	385	204	589	793	309	694	-0.5407
TOC	11.68	11.37	23.04	34.41	17.2	28.89	-0.0496
Ttl Dis Solids	846	251	1096	1347	380	1225	-0.6478
Benzene	NA	NA	NA	NA	NA	NA	
1,1-dichloroethane	NA NA	NA	NΛ	NA	NA	NA	
1,1-dichloroethene	NA NA	NA	NA	NA	NA	NA .	
Toluene	NA NA	NA NA	NA NA	NA	NA	NA	
1,1,1-trichloroethane	NA	NA	NA	NA	NA	NA	

All values in mg/L unless noted otherwise

NA All sampling events below detection limit

<sup>\* 95%</sup> CI computed for all parameter except pH. A two sided 99% CI was computed.

# Dixon Marquette Cement Company Monitoring Well 6S

Parameter	MW6-S	MW6-S	MW6-S	MW6-S	MW6-S	MW6-S	MW6-S
	26-Jul-99	09-Sep-99	01-Nov-99	02-Mar-00	23-Jun-09	29-Sep-00	29-Nov-00
Aluminum	3.01	0.08	0.116	0.166	0.04	0.127	1.74
Antimony	0.001	0.001	0.001	0.001	0.001	0.001	0.001
Arsenic	0.004	0.001	0.001	0.003	0.001	0.001	0.003
Barium	0.350	0.331	0.271	0.276	0.143	0.251	0.181
Beryllium	0.001	0.001	0.001	0.001	0.001	0.001	0.001
Boron	0.510	0.630	0.688	0.431	0.05	0.292	0.364
Cadmium	0.001	0.001	0.001	0.001	0.001	0.001	0.001
Calcium	345	24.6	232	398	149	233	256
Chromium	0.007	0.001	0.002	0.005	0.002	0.003	0.003
Cobalt	0.02	0.008	0.009	0.030	0.001	0.007	0.011
Copper -	0.024	0.006	0.004	0.022	0.001	0.012	0.02
Lead	0.015	0.001	0.001	0.001	0.001	0.001	0.008
Iron	5.34	0.693	1.06	5.55	0.284	0.77	4.96
Magnesium	78	31.4	51.9	99.2	20.3	47.7	<b>5</b> 7.6
Manganese	3.11	0.694	2.44	2.13	0.094	1.61	1.79
Mercury	0.0002	0.0002	0.0002	0.0002	0.0002	0.0002	0.0002
Nickel	0.033	0.017	0.017	0.038	0.003	0.009	0.023
Potassium	64.6	70.2	68.6	59.8	31.2	49.5	39.7
Selenium	0.003	0.001	0.002	0.005	0.001	0.003	0.004
Sodium	35.1	55	27.3	33.4	53.6	64.9	55.6
Thailium	0.001	0.001	0.001	0.001	0.001	0.001	0.001
Vana <b>dium</b>	0.023	0.001	0.003	0.002	0.005	<b>0</b> .006	0.013
Zinc	0.041	0.010	0.011	0.003	0.901	0.013	0.023
Alkalinity, total	369	354	390	388	417	384	420
Alkalinity, phenoi.				1.0	0.001	1.9	1.0
Alkalinity, bicarb	360	325	350	379	410	350	387
Ammonia nitrogen	0.10	0.10	0.10	0.10	0.10	0.10	0.10
Chloride	187	142	105	133	71.1	246	186
COD	20	20	20	20	100	123	94.1
Fluoride	0.29	0.17	0.22	0.19	0.41	0.27	0.3
Nitrate as N	0.10	0.1 <b>0</b>	0.10	0.10	0.10	0.10	0.10
pH Upper Limit (stnd units)	7.4	7.9	7.8	7.8	7.1	7.1	7.8
pH Lower Limit (stnd units)							
Sulfate	214	332	314	242	106	186	254
TOC	21.1	13.2	5.59	13.7	13.2	2.01	2.53
Ttl Dis Solids	367	865	820	789	566	842	830
Beazene	0.005	0.005	0.005	0.005	0.005	0.005	0.005
1,1-dichloroethane	0.005	0.005	0.005	0.005	0.005	0.005	0.005
1,1-dichloroethene	0.005	0.005	0.005	0.005	0.005	0.005	0.005
Toluene	0.005	0.015	0.005	0.005	0.005	0.005	0.005
1,1,1-trichloroethane	0.005	0.005	0.005	0.005	0.005	0.005	0.005

All values in mg/L unless noted otherwise

XXX Non detected Concentration

XXX Concentration > Average + 1 Standard Deviation

### Dixon Marquette Cement Company Monitoring Well 6S

Parameter	Average	Stnd Dev	Average + 1 Stnd Dev	Average + 2 Stnd Dev	99.5% CI*	Ave + 99.5% CI*	Slope of Trend
	30-Mar-01	30-Mar-01	30-Mar-01	30-Mar-0!	30-Mar-01	30-Mar-01	26-Jan-01
Aluminum	0.75	1.17	1.921	3.088	1.768	2.522	-0.0011
Antimony	0.001	0.000	0.001	0.001	0.000	0.001	
Arsenic	0.002	0.001	0.003	0.005	0,002	0.004	-0.0000009
Barium	0.258	0.075	0.332	0.407	0.113	0.371	-0.0003
Beryllium	0.001	0.000	0.001	. 0.001	0.000	0.001	
Boron	0.424	0.216	0.640	0.856	0.327	0.751	-0.0008
Cadmium	0.001	0.000	0.001	0.001	0.000	0.001	
Calcium	233.9	123.0	357	480	186	420	0.0385
Chromium	0.003	0.002	0.005	0.007	0.003	0.006	-0.000002
Cobalt	0.012	0.010	0.022	0.032	0.015	0.027	-0.00001
Copper	0.013	0.009	0.022	0.031	0.014	0.027	0.000001
Lead	0.004	0.006	0.010	0.015	0.008	0.012	-6.00E-06
Iron	2.665	2.465	5.131	7.596	3.734	6.399	-0.0004
Magnesium	55.16	26.81	82.0	108.8	40.6	95.8	-0.0237
Manganese	1.695	1.027	2.722	3.749	1.555	3.251	-0.0018
Mercury	0.0002	0.0000	0.0002	0.0002	0.0000	0.0002	
Nickel	0.020	0.012	0.032	0.045	0.019	0.039	-0.00002
Potassium	55	15	69.9	84.9	22.8	77.6	-0.0643
Selenium	0.003	0.001	0.004	0.006	0.002	0.005	0.000002
Sodium	46.4	14.2	60.6	74.9	21.5	68.0	0.0046₩
Thallium	0.001	0.000	0.001	0.001	0.000	0.001	
Vanadium	0.008	0.008	0.015	0.023	0.012	0.019	-0.000004
Zinc	0.015	0.014	0.028	0.042	0.021	0.035	
Alkalinity, total	389	24	413	436	36	425	0.0394
Alkalinity, phenol.	0.8	0.5	1.2	1.7	1.7	2.4	
Alkalinity, bicarb	366	28	394	, 422	43	409	0.0792
Ammonia nitrogen	0.10	0.00	0.10	0.10	0.00	0.10	
Chloride	152.9	58.3	211	270	88	241	0.0876
COD	56.7	46.6	103	150	71	127	0.241
Fluoride	0.26	0.08	0.35	0.43	0.12	0.39	0.0002
Nitrate as N	0.10	0.00	0.10	0.10	0.00	0.10	
pH Upper Limit (stnd units)	7.6	0.4	7.9	8.3	0.5	8.1	-0.0006
pH Lower Limit (stnd units)						7.0	
Sulfate	235	77	312	389	117	352	-0.1863
тос	10.2	7.0	17.22	24.24	10.64	20.83	-0.026
Ttl Dis Solids	797	105	902	1008	160	957	-0.1683
Benzene	NA	NA	NA	NA	NA	NA	
1,1-dichloroethane	NA	NA	NA	NA	NA	NA	
1, i-dichloroethene	NA	NA	NA	NA	NA	NA	
Toluene	NA	NA	NA	NA	NA	NA	
1,1,1-trichloroethane	NΛ	NA	NA	NA	NA	NA	

All values in mg/L unless noted otherwise

NA All sampling events below detection limit

<sup>\* 95%</sup> CI computed for all parameter except pH. A two sided 99% CI was computed.

Parameter	MW7-S	MW7-S	MW7-S	MW7-S	MW7-S	MW7-S	MW7-S
	26-Jul-99	09-Sep-99	01-Nov-99	02-Mar-00	23-Jun-00	29-Sep-00	29-Nov-00
Aluminum	0.88	0.30	0.443	0.662	0.04	0.292	0.09
Antimony	0.001	0.001	0.001	0.003	0.001	0.001	0.001
Arsenic	0.028	0.025	0.044	0.017	0.007	0.013	0.026
Barium	0.045	0.047	0.046	0.055	0.033	0.044	0.025
Beryllium	0.001	0.001	0.001	0.001	0.001	0.001	0.001
Boron	0.25	0.19	0.216	1.95)	0.05	0.204	0.195
Cadmium	0.001	0.001	0.001	0.001	0.001	0.001	0.001
Calcium	23.9	14.7	11.9	43.4	40.8	29	14.2
Chromium	0.004	0.002	0.002	0.005	0.002	0.001	0.002
Cobalt	0.001	0.001	0.001	0.001	0.001	0.001	0.001
Copper	0.005	0.007	0.008	0.006	0.001	0.006	0.004
Lead ,	0.003	0.002	0.002	0.001	0.001	0.001	0.001
Iron	1.98	0.789	0.629	1.32	0.109	0.663	0.142
Magnesium	10.8	5.8	4.55	13.6	15.1	10.6	4.6
Manganese	0.154	0.133	0.085	0.157	0.002	0.103	0.024
Mercury	0.0002	0.0002	0.0002	0.0002	0.0002	0.0002	0.0002
Nickel	0.012	0.014	0.014	0.007	0.002	0.011	0.009
Potassium	656	717	793	503	406	626	727
Selenium	0.001	0.002	0.002	0.005	0.001	0.003	0.002
Sodium	37.6	42.9	49.9	37.4	25.8	44.2	48.8
Thallium	0.001	0.001	0.001	0.001	0.001	0.001	0.001
Vanadium	0.016	0.012	0.022	0.015	0.007	0.013	0.006
Zinc	0.027	0.015	0.016	0.027	0.001	0.023	0.001
Alkalinity, total	396	466	534	348	341	384	486
Alkalinity, phenol.		·		41	1.0	92	160
Alkalinity, bicarb	285	265	475	154	333	254	297
Ammonia nitrogen	0.10	0.10	0.10	0.10	0.10	0.94	0.10
Chloride	70.1	90.2	97.7	56.3	48.8	105	110
COD	20	20	20	20	10.5	34.3	20
Fluoride	1.64	1.95	2.43	0.8	0.65	1.3	1.84
Nitrate as N	0.10	0.10	0.10	0.10	0.10	0.10	0.10
pH Upper Limit (stnd units)	10	10.2	9.8	9.4	8.9	9.4	10.2
pH Lower Limit (stnd units)							
Sulfate	401	475	494	345	336	448	474
TOC	17.2	10.7	15.2	14.4	18	7	11.3
Ttl Dis Solids	1047	1294	1423	901	802	1156	1308
Benzene	0.005	0.005	0.005	0.005	0.005	0.005	0.005
1,1-dichloroethane	0.005	0.005	0.0389	0.013	0.0252	0.0639	0.0581
1,1-dichloroethene	0.005	0.0061	0.005	0.005	0.005	0.0061	0.0082
Toluene	0.005	0.005	0.005	0.005	0.005	0.005	0.005
1,1,1-trichloroethane	0.006	0.005	0.005	0.005	0.0057	0.005	0.005

All values in mg/L unless noted otherwise

XXX Non detected Concentration

XXX Concentration > Average + 1 Standard Deviation

## Dixon Marquette Cement Company Monitoring Well 7S

Parameter	Average	Stnd Dev	Average + 1 Stnd Dev	Average + 2 Stnd Dev	99.5% CI*	Ave + 99.5% CI*	Slope of Trend
	30-Mar-01	30-Mar-01	30-Mar-01	30-Mar-01	30-Mar-01	30-Mar-01	26-Jan-01
Aluminum	0.387	0.302	0.689	0.991	0.457	0.844	-0.0011
Antimony	0.001	0.001	0.002	0.003	0.001	0.002	-0.0000001
Arsenic	0.023	0.012	0.035	0.047	0.018	0.041	-0.00003
Barium	0.042	0.010	0.052	0.062	0.015	0.057	-0.00003
Beryllium	0.001	0.000	0.001	0.001	0.000	0.001	
Boron	0.436	0.670	1.107	1.777	1.015	1.452	-0.0002
Cadmium	0.001	0.000	0.001	0.001	0.000	0.001	
Calcium	25	13	38	51	20	45	0.0147
Chromium	0.003	0.001	0.004	0.005	0.002	0.005	-0.000063
Cobalt	0.001	0.000	0.001	0.001	0.000	0.001	
Copper	0.005	0.002	0.008	0.010	0.003	0.009	-0.000006
Lead	0.002	0.001	0.002	0.003	100.0	0.003	-0.000003
Iron	0.805	0.661	1.466	2.126	1.001	1.806	-0.0023
Magnesium	9.3	4.3	13.6	18.0	6.6	15.9	0.0023
Manganese	0.094	0.061	0.155	0.217	0.093	0.187	-0.002
Mercury	0.0002	0.0000	0.0002	0.0002	0.0000	0.0002	
Nickel	0.010	0.004	0.014	0.018	0.007	0.016	-0.00001
Potassium	632.6	135.7	768.3	904.0	205.5	838.1	-0.1819
Selenium	0.002	0.001	0.004	0.005	0.002	0.004	0.000001
Sodium	40.9	8.3	49.2	57.5	12.5	53.5	0.0018
Thallium	0.001	0.000	0.001	0.001	0.000	0.001	
Vanadium	C 013	0.005	0.018	0.024	0.008	0.021	-0.00002
Zinc	0.016	0.011	0.027	0.038	0.017	0.033	-0.00003
Alkalinity, total	422.14	73.86	496.00	569.87	111.87	534.01	0.0687
Alkalinity, phenol.	7~.5	69	142.1	210.8	231	305	0.453
Alkalinity, bicarb	294.71	96.96	391.68	488.64	146.86	441.57	-0.081
Ammonia nitrogen	0.2	0.3	0.5	0.9	0.5	0.7	0.0007
Chloride	82.59	24.25	106.84	131.09	36.73	119.32	0.0354
COD	20.69	6.97	27.66	34.62	10.56	31.24	0.0084
Fluoride	1.52	0.64	2.16	2.79	6.97	2.48	-0.0012
Nitrate as N	0.10	0.00	0.10	0.10	0.0	0.10	
pH Upper Limit (stnd units)	9.7	0.5	10.2	10.7	0.7	10.4	-0.0008
pH Lower Limit (stnd units)						9.0	
Sulfate	425	65	489	554	98	523	-0.0167
TOC	13.4	3.9	17.3	21.3	5.9	19.3	-0.0088
Ttl Dis Solids	1133	228	1361	1589	345	1478	-0.1319
Benzene	NA	NA	NA	NA	NA	NA	
1,1-dichlor oethane	0.0299	0.0244	0.054	0.0787	0.037	0.0669	0.1027
1,1-dichleroethene	0.0058	0.0012	0.007	0.0082	0.002	0.0076	0.0038
Toluene	NA	NA	NA	NA	NA	NΛ	
1,1,1-trichloroethane	0.0052	0.0004	0.006	0.0061	0.001	0.0059	-0.0007

All values in mg/L unless noted otherwise

NA All sampling events below detection limit

<sup>\* 95%</sup> Ci computed for all parameter except pH. A two sided 99% CI was computed.

#### Dixon Marquette Cement Company Monitoring Well 8S

Parameter	MW8-S	MW8-S	MW8-S	MW8-S	MW8-S	MW8-S	MW8-S
•	26-Jul-99	09-Sep <b>-99</b>	01-Nov-99	02-Mar-00	23-Jun-00	29-Sep-00	29-Nov-00
Aluminum	0.51	0.29	0.665	0.435	0.04	0.109	0,233
Antimony	0.001	0.001	0.001	0.001	0.001	0.001	0.001
Arsenic	0.001	0.001	100.0	0.005	0.001	0.001	0.001
Barium	0.073	0.066	0.056	0.172	0.044	0.055	0.057
Beryllium	0.001	0.001	6.001	0.002	0.001	0.001	0.001
Boron	0.25	0.25	0.225	0.253	0.05	0.199	0.232
Cadmium	0.001	0.001	0.001	0.003	0.001	0.001	0.001
Calcium	265	226	223	268	224	234	222
Chromium	0.004	0.003	0.006	0.04	0.007	0.005	0.006
Cobalt	0.002	0.002	0.002	0.035	0.001	0.001	0.001
Copper	0.015	0.014	0.011	0.113	0.001	0.007	0.007
Lead	0.003	0.004	0.003	0.004	0.001	0.001	0.002
Iron	1.6	1.19	1.57	5.3	0.415	0.549	0.953
Magnesium	110	94.4	94.3	125	92.4	93.3	91.1
Manganese	<b>0.363</b>	0.233	0.255	4.3	<b>0</b> .001	0.024	0.167
Mercury	0.0002	0.0002	0.0002	0.0002	0.0002	9.0002	0.0002
Nickel	0.010	0.010	0.008	0.058	0.002	0.003	0.005
Potassium	265	263	265	217	174	254	280
Selenium	0.003	0.003	0.003	0.014	0.001	0.002	0.006
Sodium	8.2	39.6	40.1	43.1	43.4	52.4	51.4
Thallium	0.001	9.001	0.001	0.002	0.001	0.005	0.001
Vanadium	0.003	105.0	0.003	0.001	0.004	0.003	0.003
Zinc	0.044	0.027	ባ.042	0.363	0.001	0.025	0.021
Alkalinity, total	271	270	314	284	316	274	278
Alkalinity, phenol.				! 0	1.0	1.0	1.0
Alkalinity, blearb	265	265	300	271	314	262	257
Ammonia nitrogen	0.10	0.10	0.10	0.10	0.10	0.10	0.10
Chloride	119	121	113	162	181	179	198
COD	20	20	20	30.3	158 J	30	20
Fluoride	0.26	0.16	0.19	0.14	0.26	0.22	0.21
Nitrate as N	2.73	2.47	2.68	1.68	2.95	3.01	1.86
pH Upper Limit (stnd units) pH Lower Limit (stnd units)	7.8	8.3	7.8	7.1	7.2	7.3	7.9
Sulfate	818	818	753	85/2	686	0.47	978
TOC	9.8	5.86	9.8	14.4	15.1	847	865
Til Dis Solids	1135	1190	7.8 1191			2.05	2.5
1 (1 1715 70HU3	1133	1190	1191	1161	1125	1200	1252
Benzene	0.005	0.005	0.005	0.005	0.005	6 005	0.005
1,1-dichloroethane	0.005	0.005	0.005	0.005	0.005	0.005	0.005
1,1-dichloroethene	0.005	0.005	0.005	0.005	0.005	0.005	0.005
Toluene	0.005	0.005	0.005	0.005	0.005	0.005	0.005
1,1,1-trichloroethane	0.005	0.005	0.005	0.005	0.005	0.005	0.005

All values in mg/L unless noted otherwise

XXX Non detected Concentration

XXX Concentration > Average + 1 Standard Deviation

Parameter	Average	Stnd Dev	Average + 1 Stnd Dev	Average + 2 Stnd Dev	99.5% CI*	Ave + 99.5% CI*	Slope of Trend
	30-Mar-01	30-Mar-01	30-Mar-01	30-Mar-01	30-Mar-01	30-Mar-01	26-Jan-01
Aluminum	0.326	0.223	0.549	0.773	0.338	0.664	-0.0008
Antimony	0.001	0 000	0.001	0.001	0.000	0.001	
Arsenic	0.002	0.002	0.603	0.005	0.002	0.004	-0.0000002
Barium	0.075	0.044	0.119	0.162	0.066	0.141	-0.00004
Beryllium	0.301	0.000	0.002	0.002	0.001	0.002	-5E-08
Boron	0.208	0.072	0.281	0.353	0.110	0.318	-0.0001
Cadmium	0.001	0.001	0.002	0.003	0.001	0.002	-0.0000001
Calcium	237	20	258	278	31	268	-0.039
Chromium	0.010	0.013	0.023	0.037	0.020	0.030	0.000002
Cobalt	0.006	0.013	0.019	0.032	0.019	0.025	-0.000004
Copper	0.024	0.040	0.064	0.103	0.060	0.084	-0.00002
Lead	0.003	0.001	0.004	0.005	0.002	0.004	-0.000005
Iron	1.654	1.672	3.325	4.997	2.532	4.186	-0.0021
Magnesium	100.1	12.7	112.8	125.5	19.3	119.3	0.0233
Manganese	0.763	1.565	2.328	3.893	2.370	3.133	0.0007
Mercury	0.0002	0.000	0.0002	0.0002	0.0000	0.0002	
Nickel	0.014	0.020	0.034	0.053	0.030	0.044	-0.00002
Potassium	245	37	283	320	56	302	-0.0296
Selenium	0.005	0.004	0.009	013	0.007	0.011	0.000001
Sodium	39.7	14.8	54.6	69.4	22.4	62.2	0.0587
Thallium	0 002	<b>j.001</b>	0.003	0.005	0.002	0.004	0.000004
Vanadium	0.003	0.001	0.004	0.005	0.002	0.004	0.000002
Zinc	0.075	0.128	0.203	0.331	0.194	0.268	-0.00006
Alkalinity, total	287	20	307	326	30	317	0.0058
Aikalinity, phenol.	1.0	0	1.0	1.0	0.0	1.0	
Alkalinity, bicarb	276	22	298	320	33	309	-0.0109
Ammonia nitrogen	01.0	0.00	0.10	0.10	0.00	0.10	
Chloride	153	35	188	223	53	206	0.173]
COD	41	52	93	144	78	119	0.0621
Fluoride	0.21	0.05	0.25	0.30	0.07	0.28	0.00004
Nitrate as N	2.48	0.52	3.00	3.52	0.8	3.27	0.004
pH Upper Limit (stnd units)	7.6	0.4	8.1	8.5	0.7	8.3	-0.001
pH Lower Limit (stnd units)						7.0	
Sulfate	803	62	865	927	94	897	0.0585
TOC	8.5	5.3	13.8	19.0	8.0	16.5	-0.0099
Til Dis Solids	1179	43	1222	1266	65	1245	0.1 <b>091</b>
Benzene	NA	NA	NA	NA	NA	NA	
1,1-dichloroethane	NA	NA	NA	NA	NA	NA	
1,1-dichloroethene	NA	NA	NA	NA	NA	NA	
Toluene	NA	NA	NA	NA	NA	NA	
1,1,1-trichloroethane	NA	NΛ	NA	NA	NA	NA	

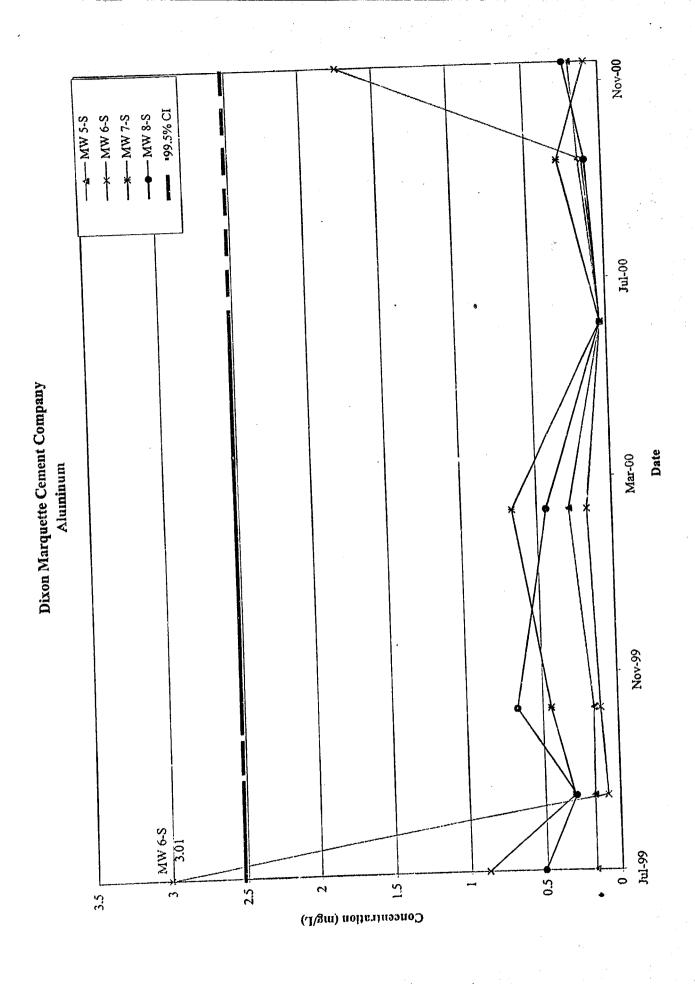
All values in mg/L unless noted otherwise

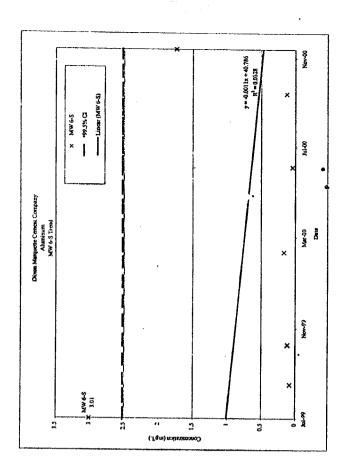
NA All sampling events below detection limit

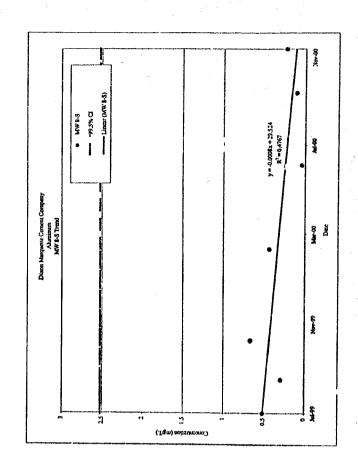
<sup>\* 95%</sup> Cl computed for all parameter except pH. A two sided 99% Cl was computed.

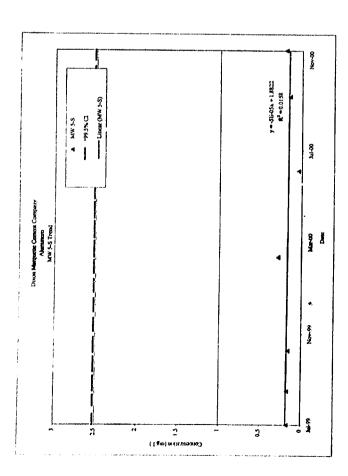
Adjusted Groundwater Standards
Dixon Marquette Cement Company -- Dixon, Illinois
Prepared by Preston Engineering, Inc.
February 27, 2001

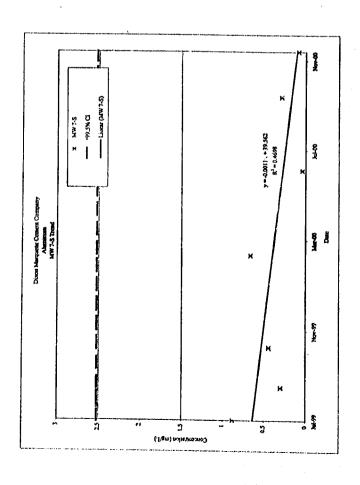
**GRAPHS** 



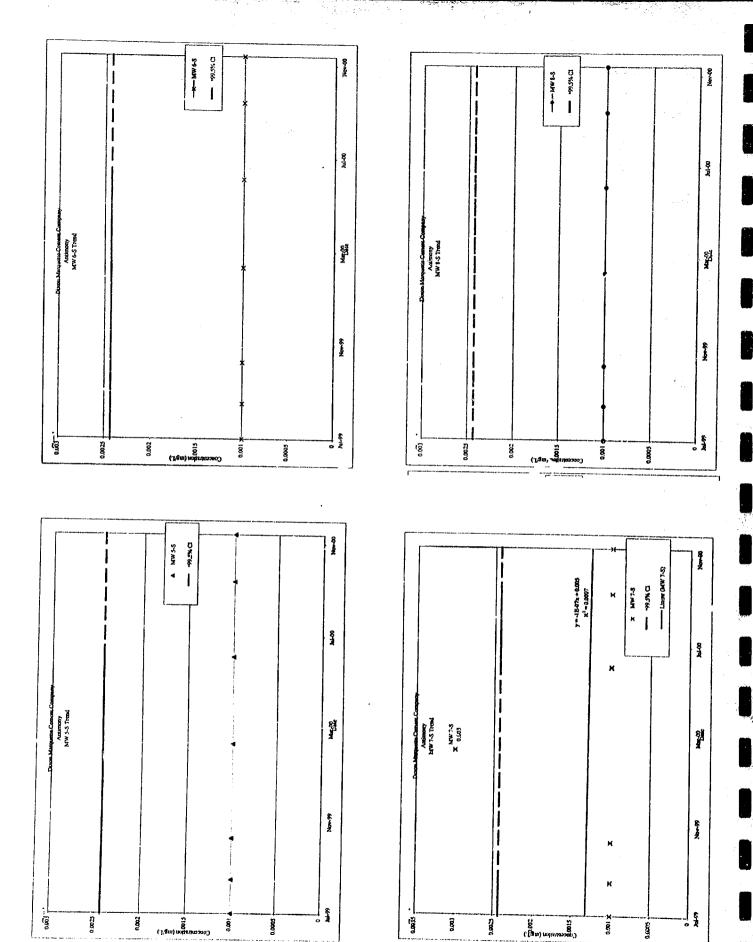




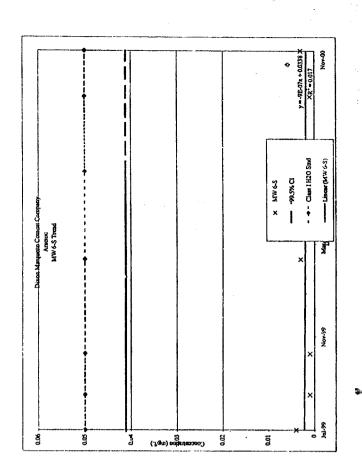


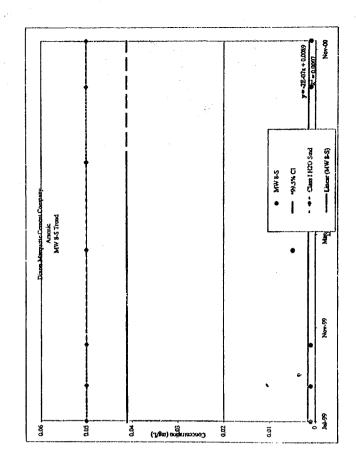


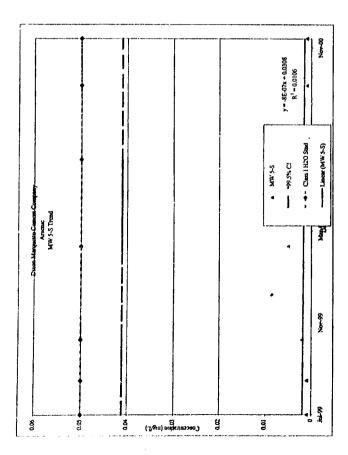
Nov-00 \*99.5% CI --- MW 8-S Jul-00 Dixon Marquette Cement Company MW 7-S Date Mar-00 Antimony Nov-99 Class I H2O Standard: 0.006 mg/L Jul-99 0.0025 0.0035 0.0015 0.003 0.002 0.0005 0.001 Concentration (mg/L)

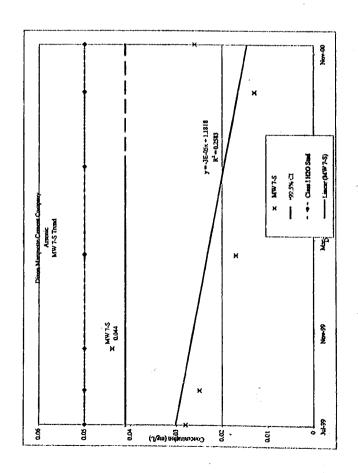


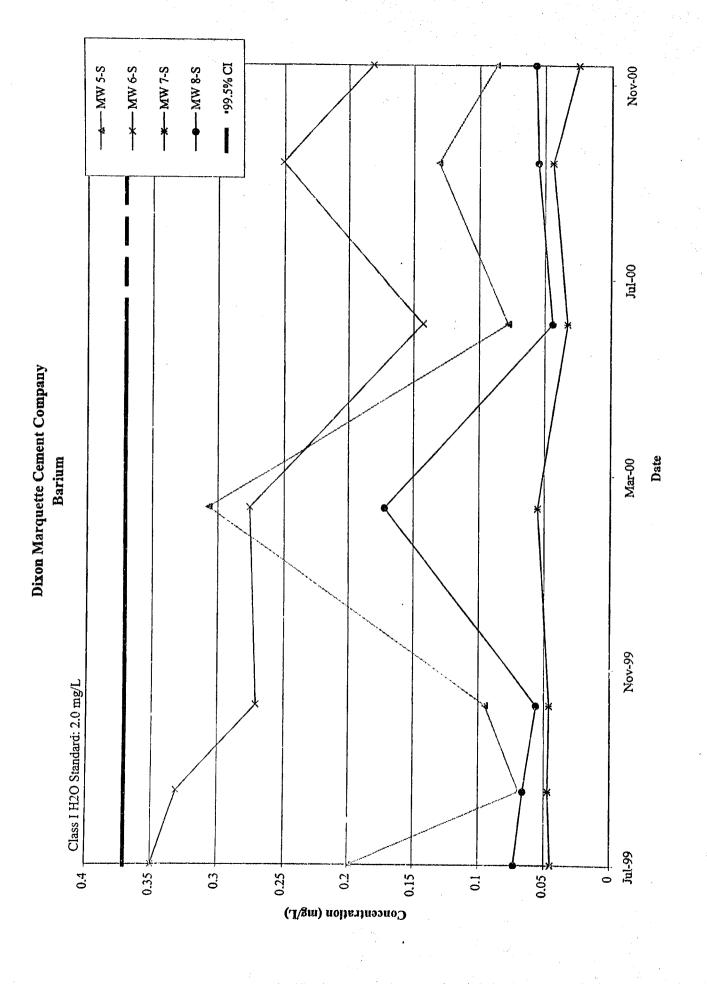
Nov-00 Jul-00 Dixon Marquette Cement Company Date Mar-00 Arsenic MW 7-S 0.044 Nov-99 ◆ Class I H2O Smd -99.5% CI \*-- MW 7-S ●—MW 8-S \*- MW 5-S Jul-99 0.06 0.05 0.04 0.03 0.02 0.01 Concentration (mg/L)

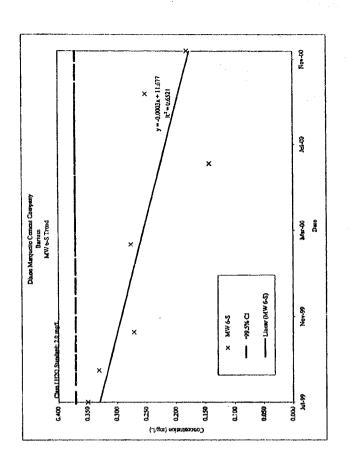


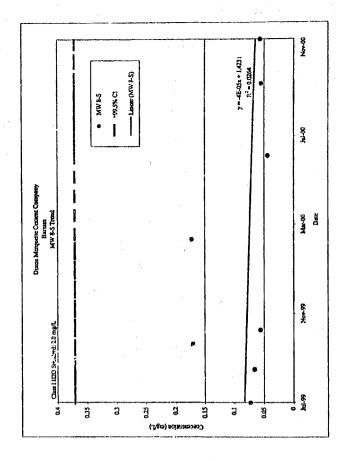


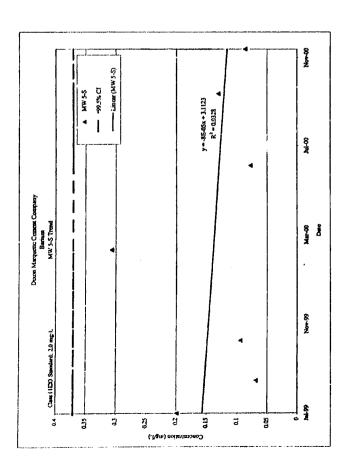


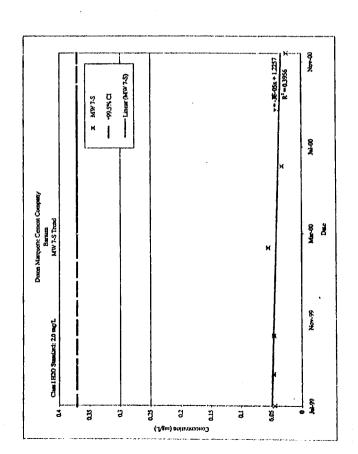




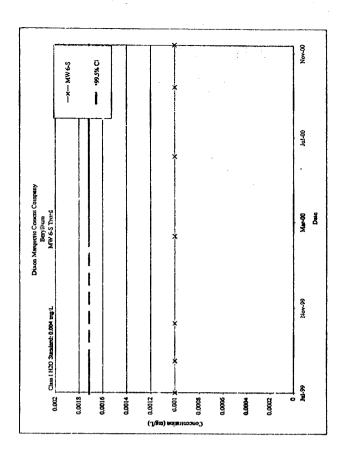


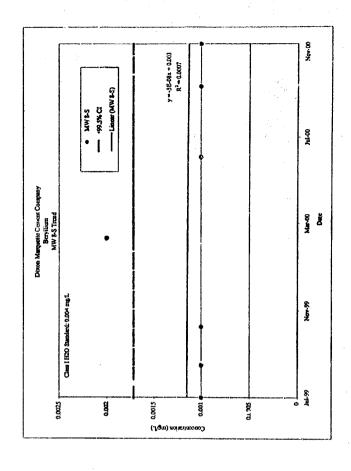


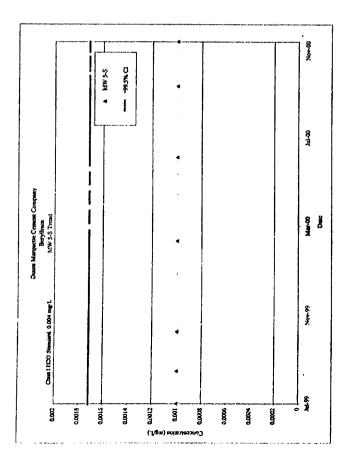


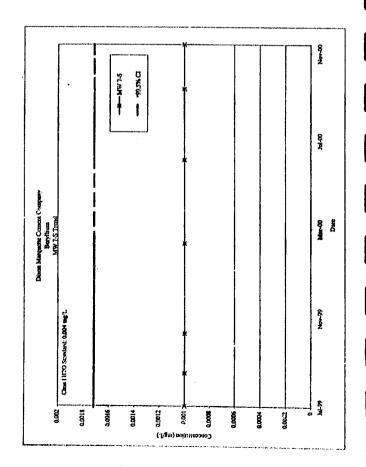


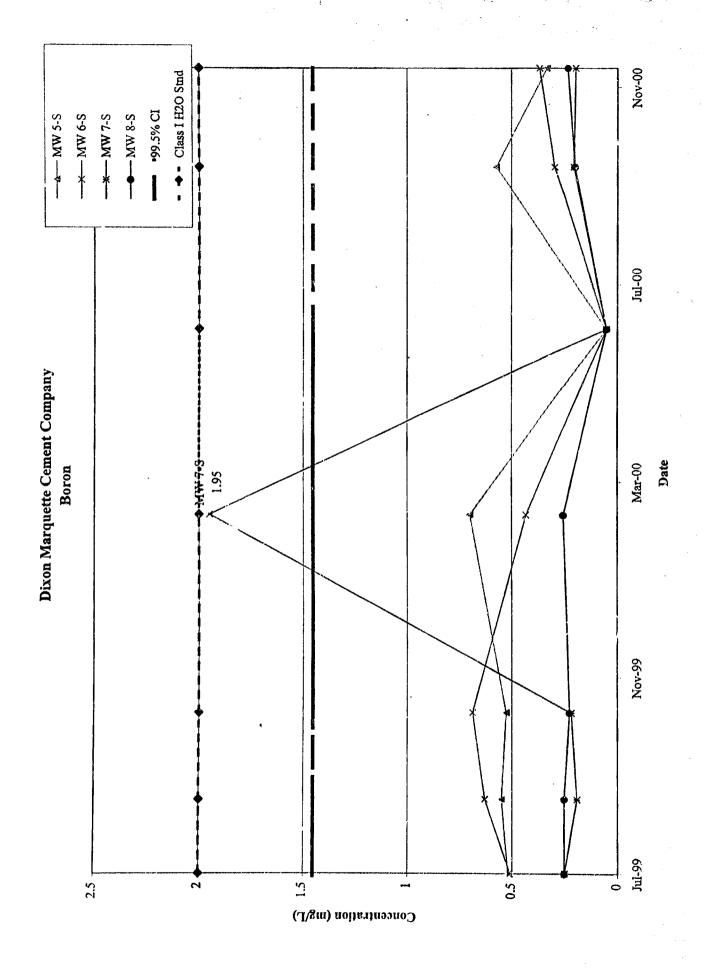
Nov-00 •99.5% CI ×-- MW 6-S -\*- MW 5-S -\*-- MW 7-S Jul-00 Dixon Marquette Cement Company Date Mar-00 Beryllium Nov-99 Class I H2O Standard: 0.004 mg/L Jul-99 0.0025 0.007 0.0015 0.0005 0.001 Concentration (mg/L)

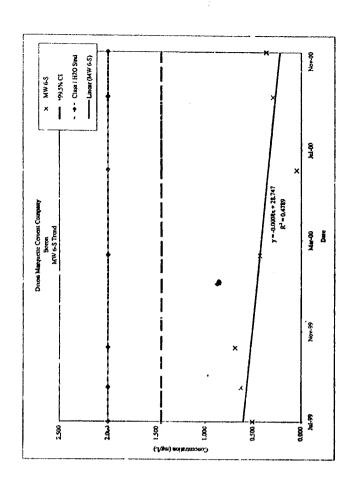


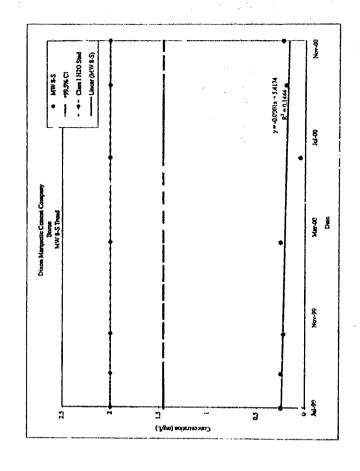


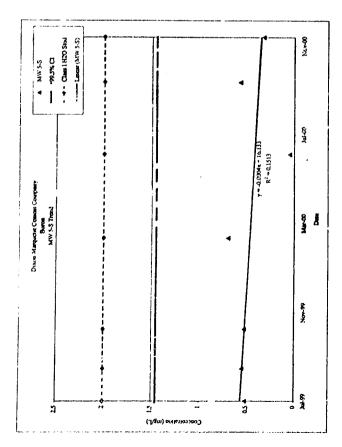


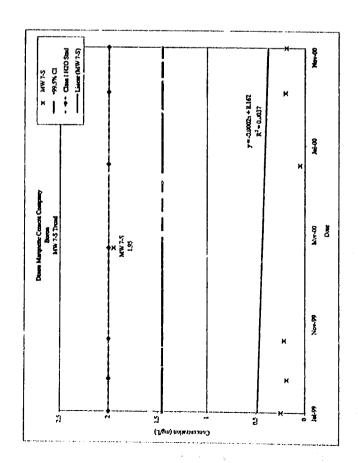






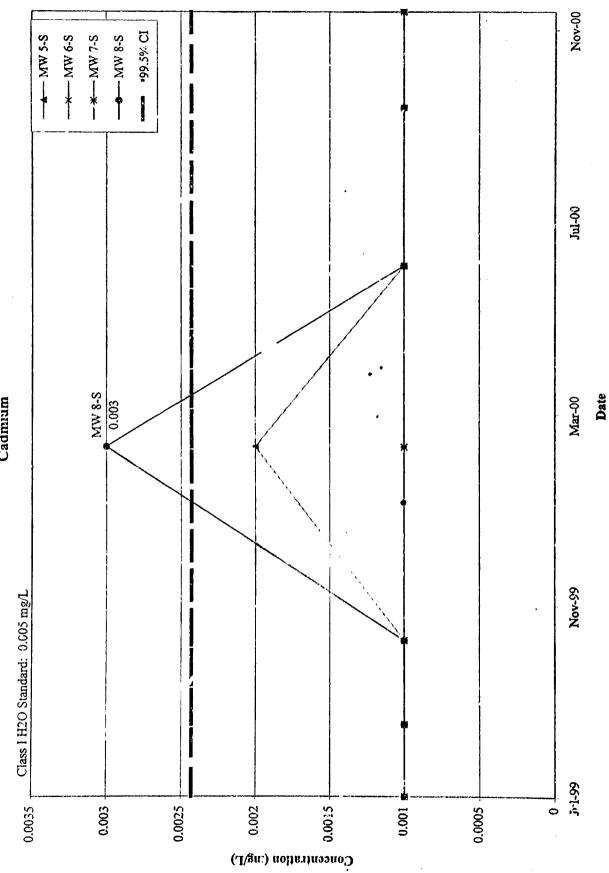


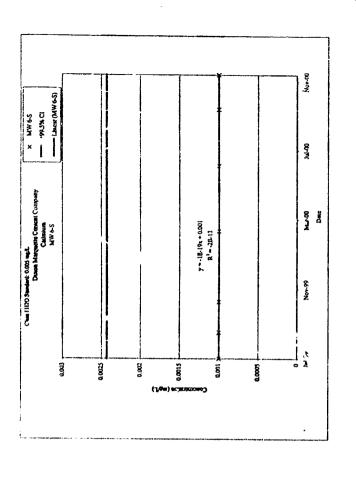


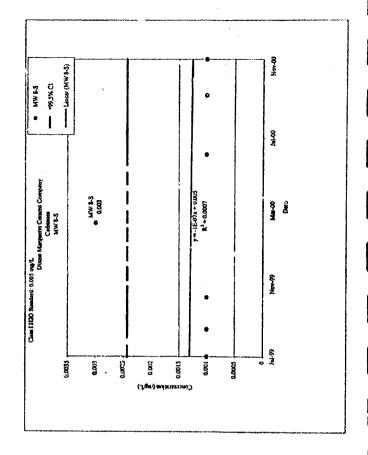


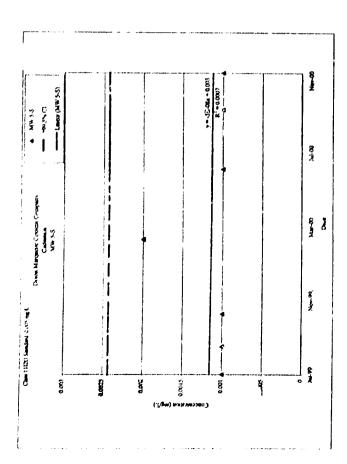
Dixon Marquette Cement Company Cadmium

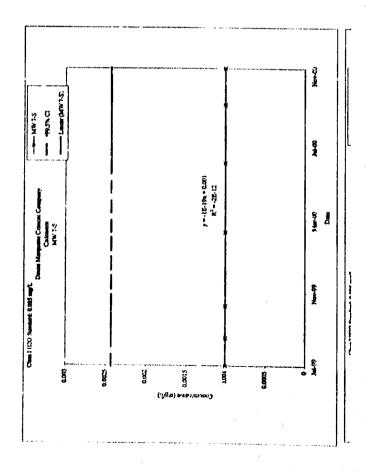
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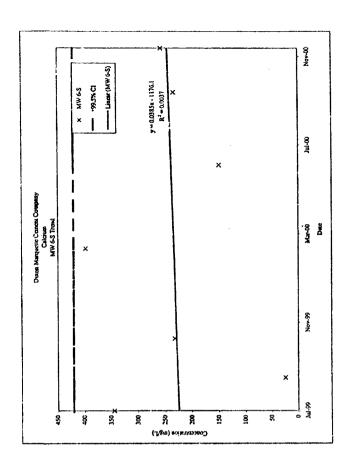


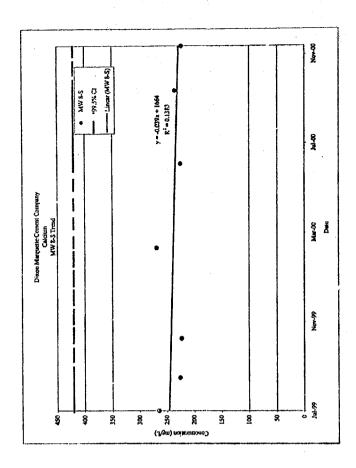


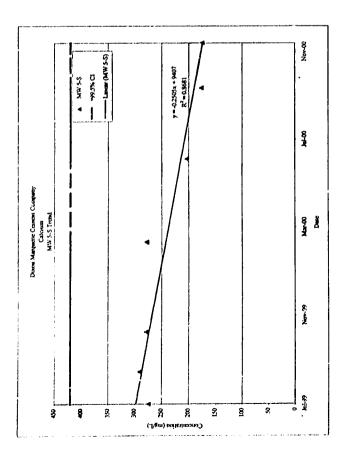


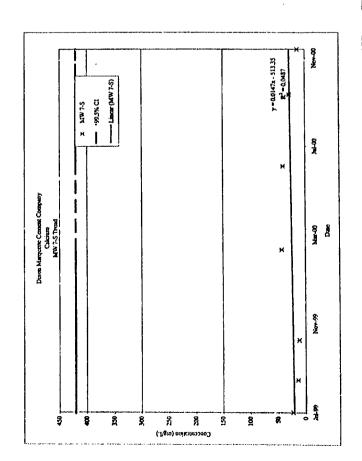


Nov-00 \*99.5% CI -\*-- WW 7-S \* WW 5-S S-9 MM −×−  $J_{\rm ul}$ -00 Dixon Marquette Cement Company Calcium Mar-00 Date Nov-99 Jul-99 450 400 350 300 250 200 150 100 20 Concentration (mg/L)



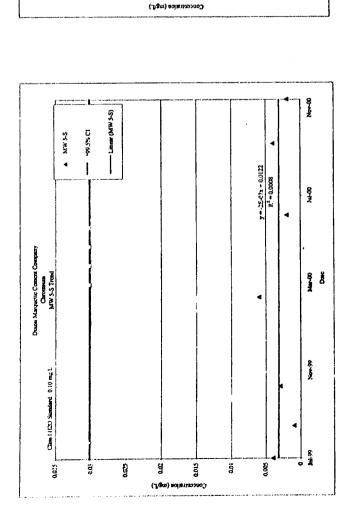






Nov-00 \*99.5% CI S-9 MW ----\*- MW 5-S --\*-- MW 7-S Jul-00 Dixon Marquette Cement Company Chromium MW 8-S Date Mar-00 Nov-99 Class I H2O Standard: 0.10 mg/L Jul-99 0.005 0 0.045 0.0 40 0.035 0.03 0.025 0.015 0.02 0.01 Concentration (mg/L)

F



Lincar (MW 6-S)

8

0.015

9.0

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5000

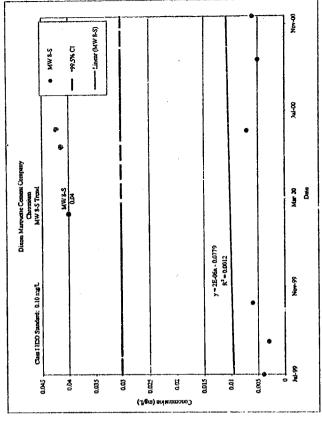
× MW 6-S

Dixon Marquetts Cement Company
Chromium
MW 6-S Trend

Class 11120 Sundard: 0.10 mg/L

0.035

90



W. \*\*\*

Nov-00

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Mar-00 Date

8-62

6-jg

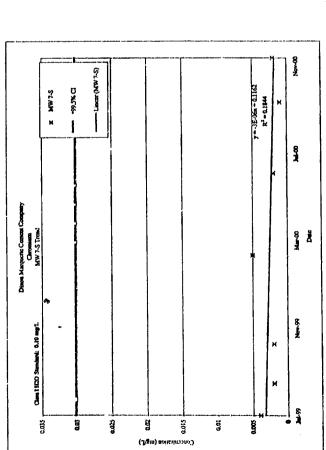
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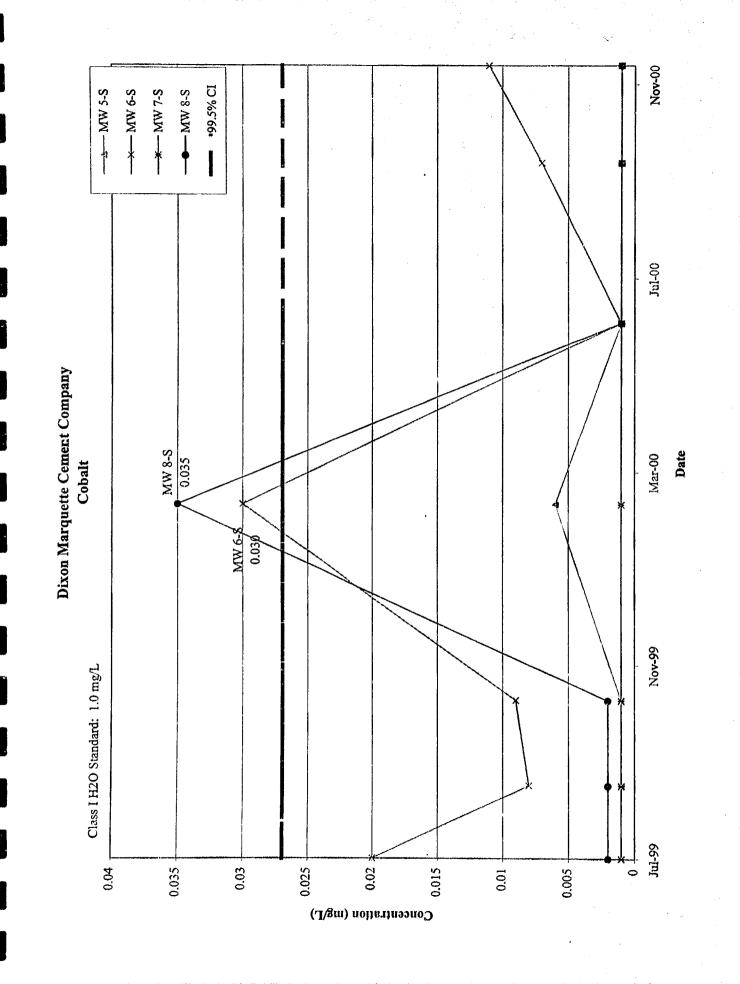
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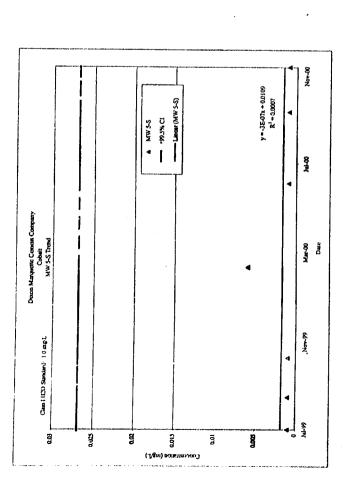
y = -2E-Dfx + 0.092 R² = 0.0518

×

×







\_\_\_\_\_Lincar (MW 6-5)

× MW6-5

Dixtu Marquette Coment Company Cobalt MW 6-5 Treed

Contract the same to see

0.035

MCW 6-S ×

g

San

0.02

0.015

Concentation (mg/L)

9

0.005

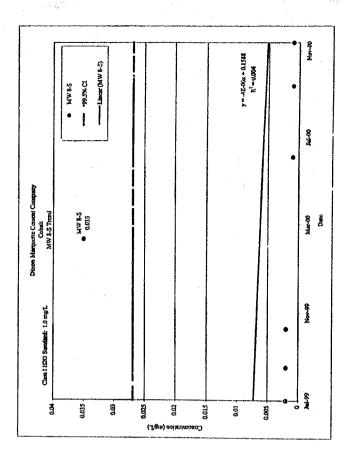
y=.1E-05x+0.5486 R<sup>2</sup>-0.0861 Nov-00

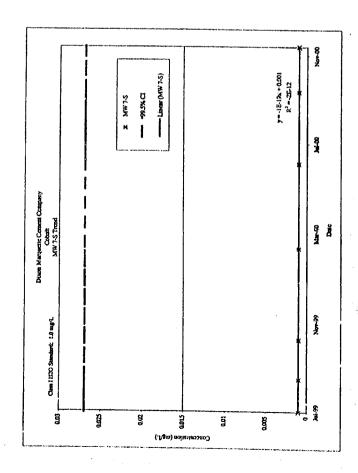
, 00-m

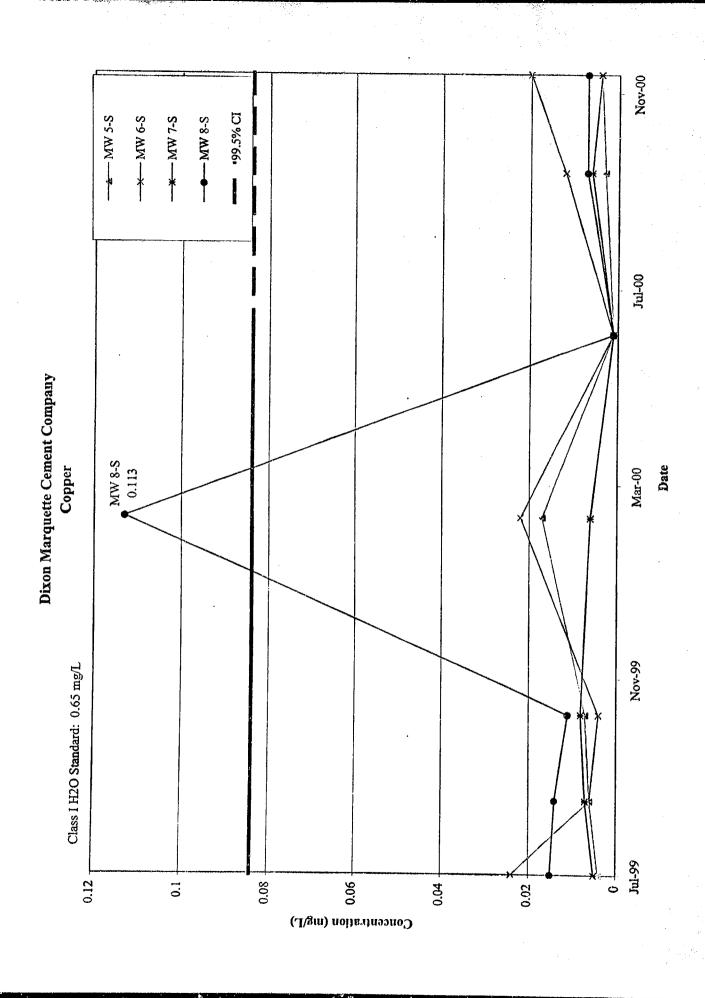
Mar-00 Date

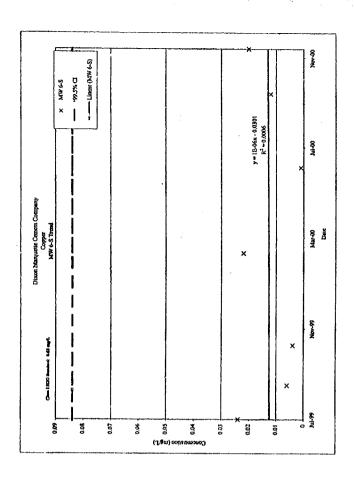
Not-99

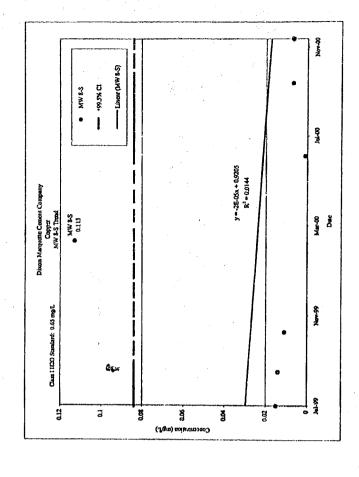
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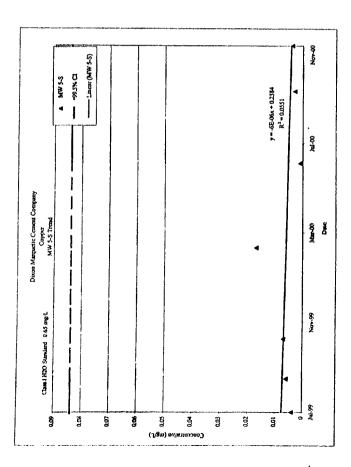


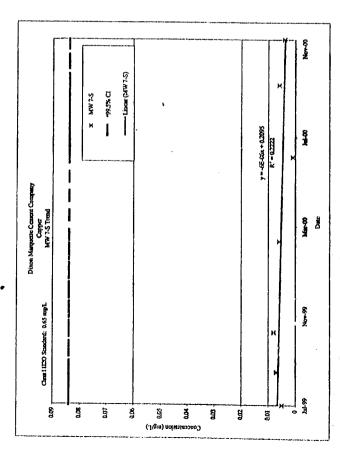


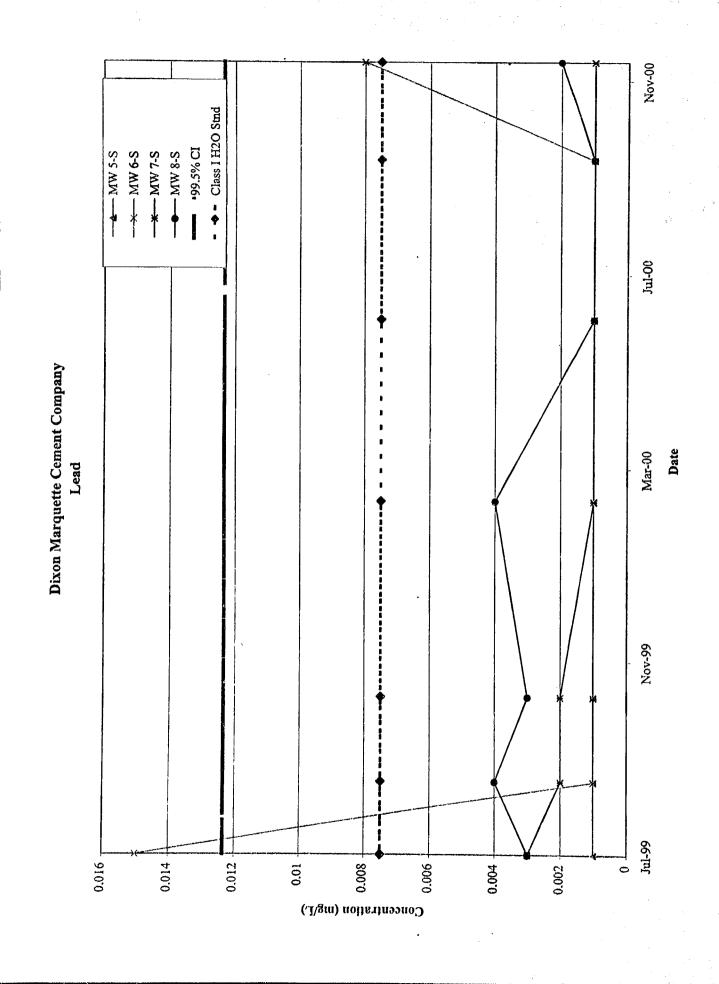


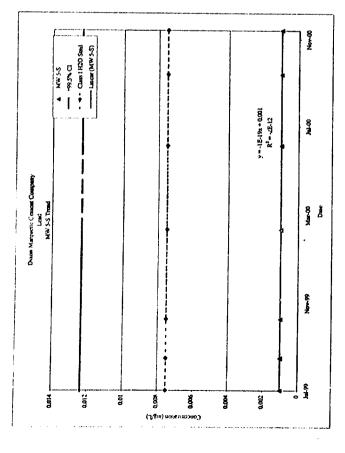


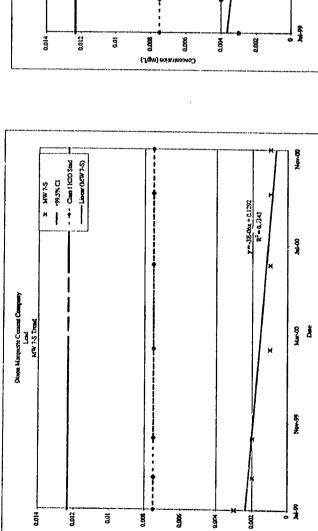




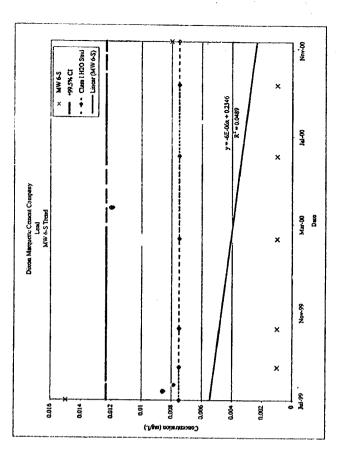


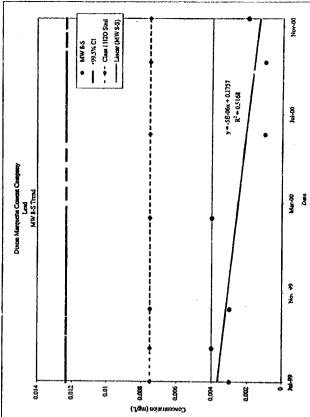




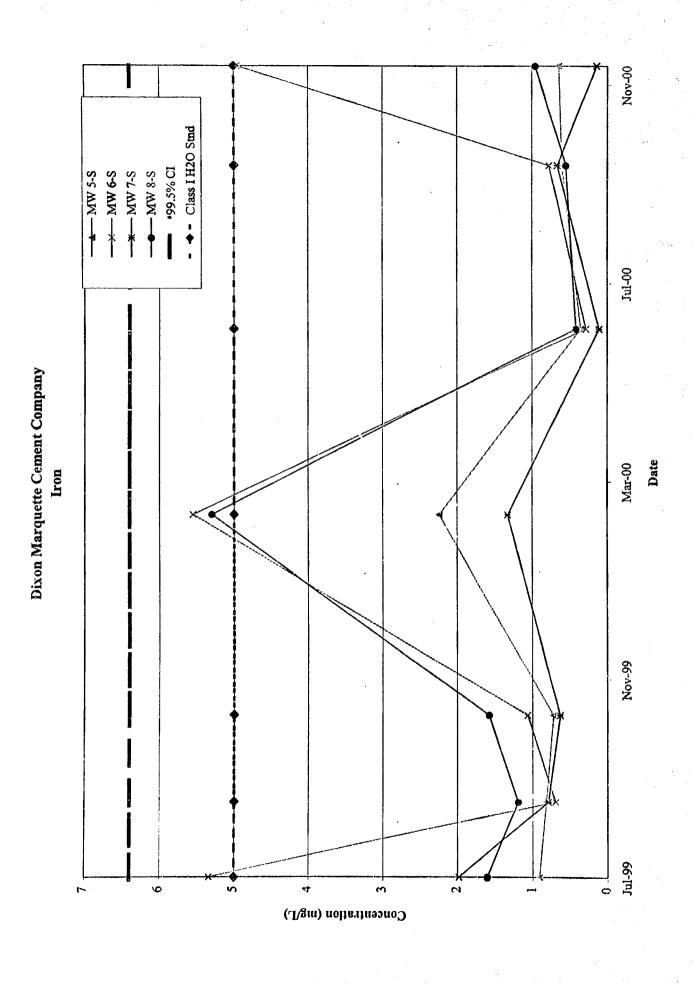


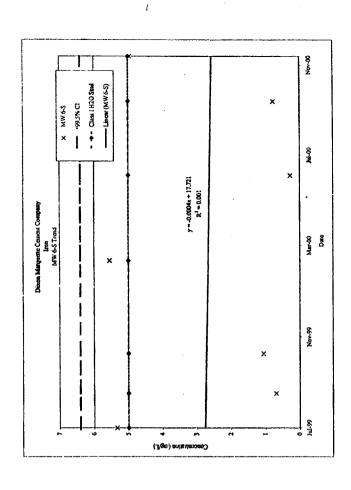
Concentiation (mg/L)

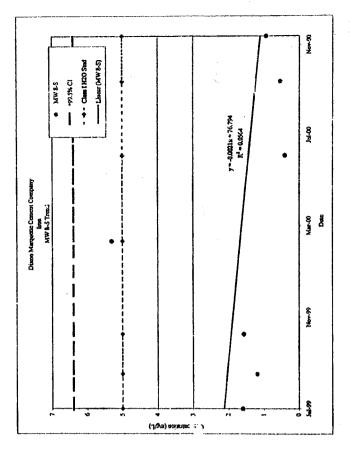


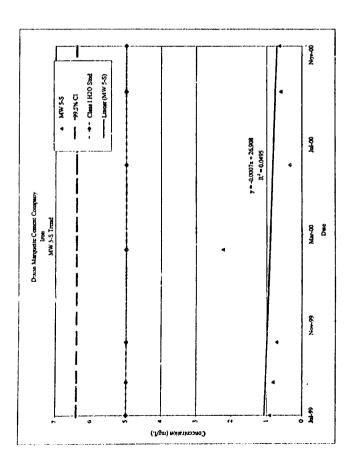


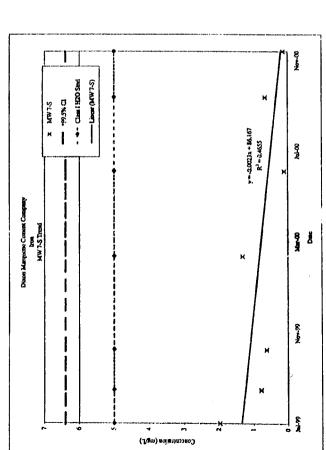
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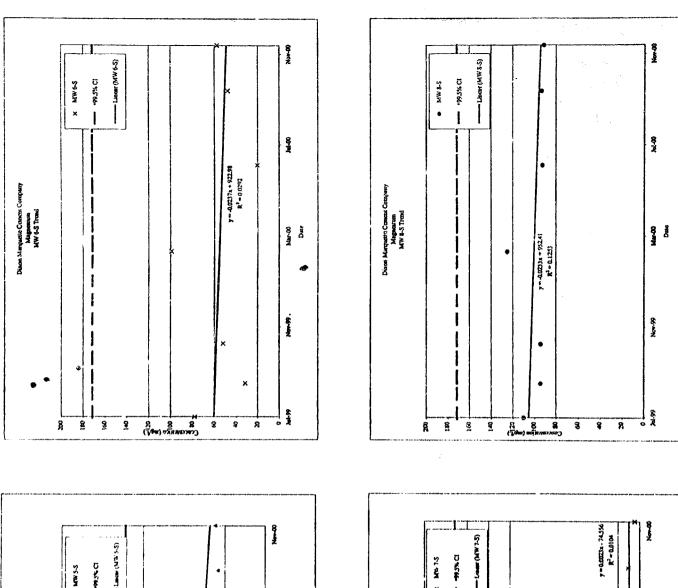


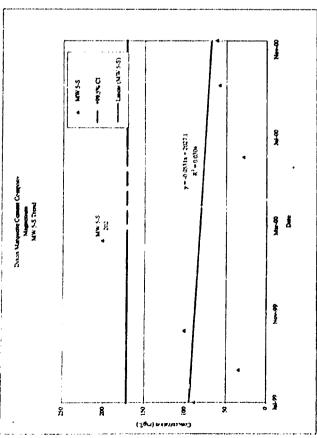




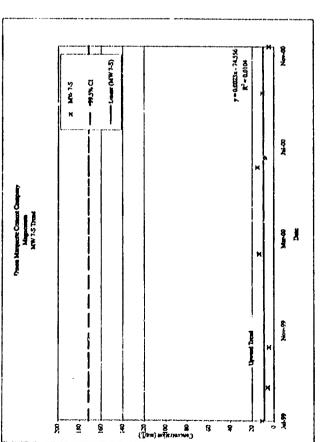


Nov-00 -99.5% CI ---\*--- MW 5-S S-9 MM ---\*-- WW 7-S --- MW 8-S Jul-00 Dixon Marquette Cement Company Magnesium MW 5-S Mar-00 Date Nov-99 Jul-99 200 250 150 50 Concentration (mg/L)

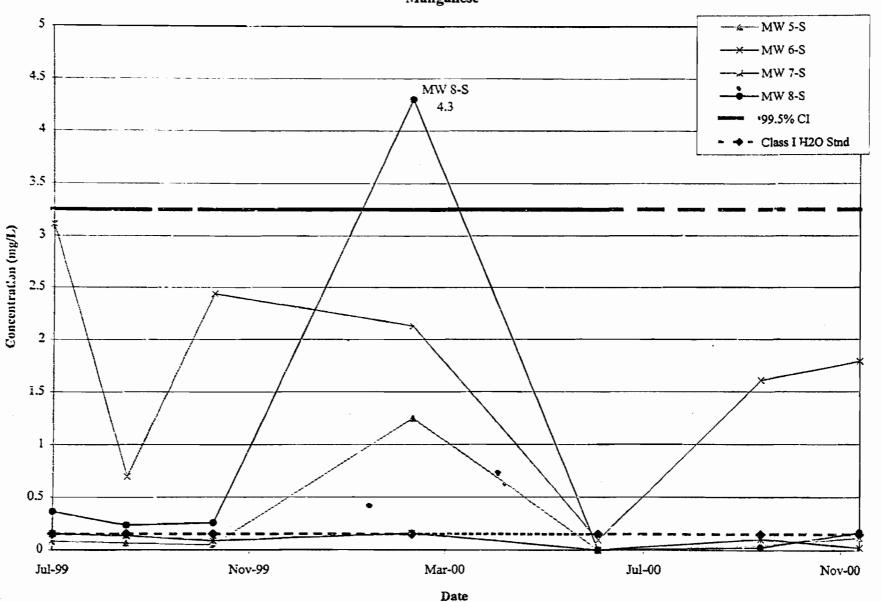


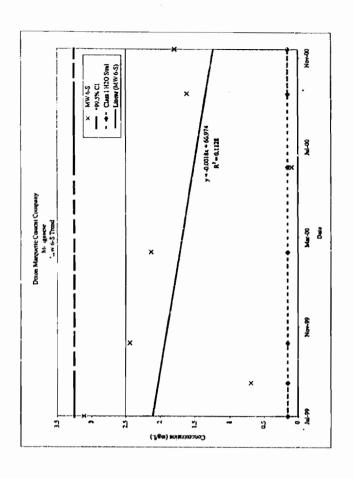


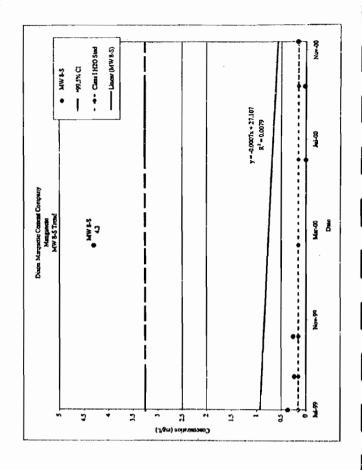
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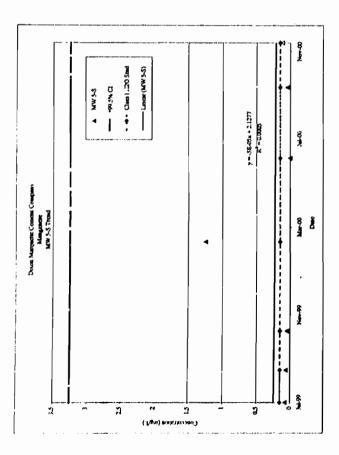


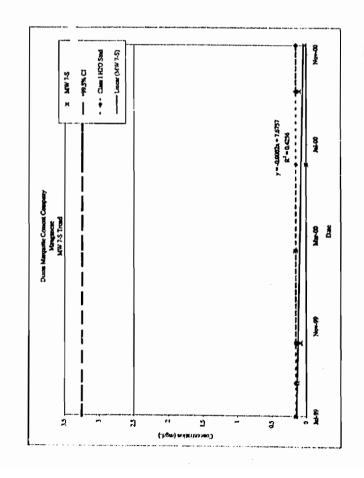
# Dixon Marquette Cement Company Manganese



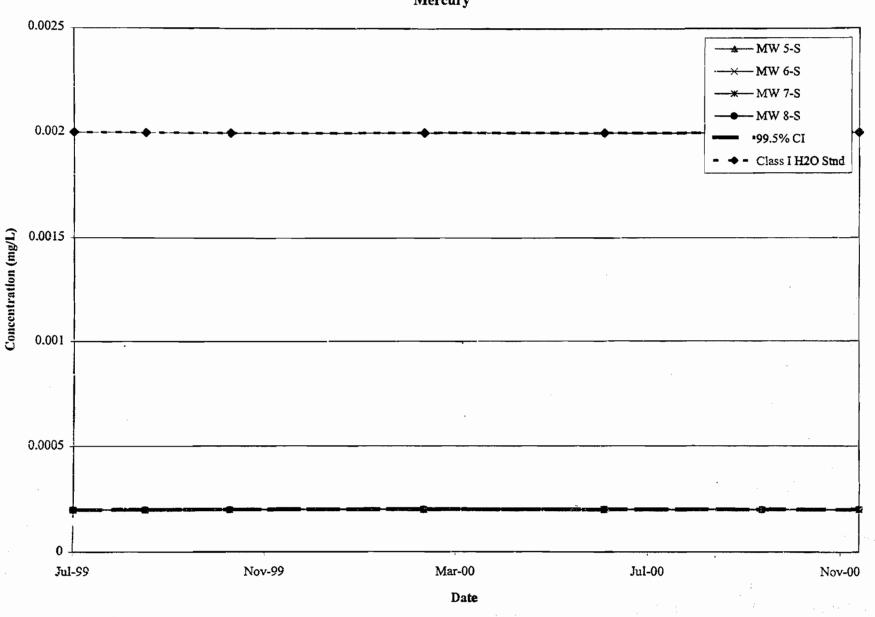




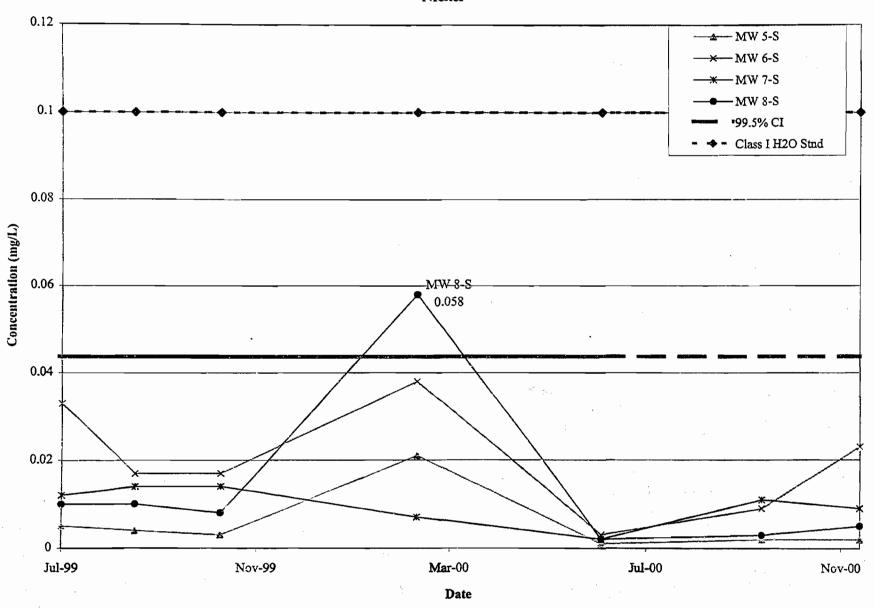


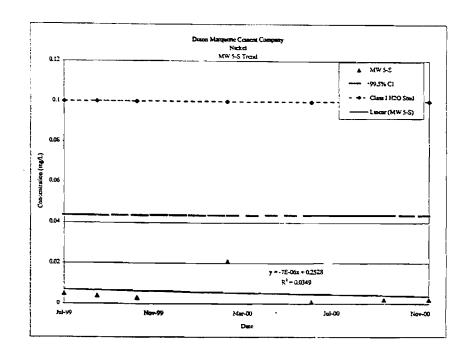


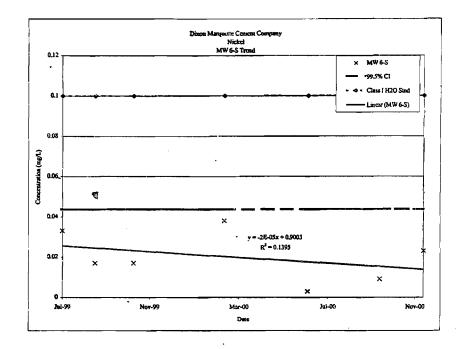
# Dixon Marquette Cement Company Mercury

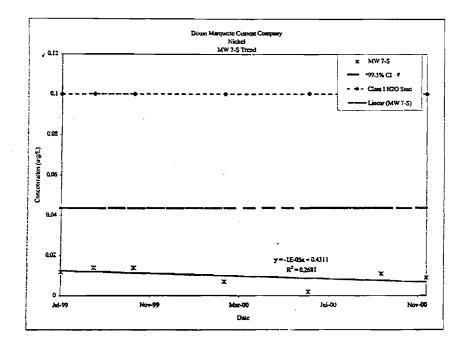


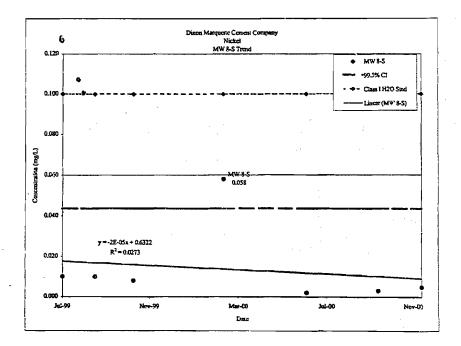
### Dixon Marquette Cement Company Nickel



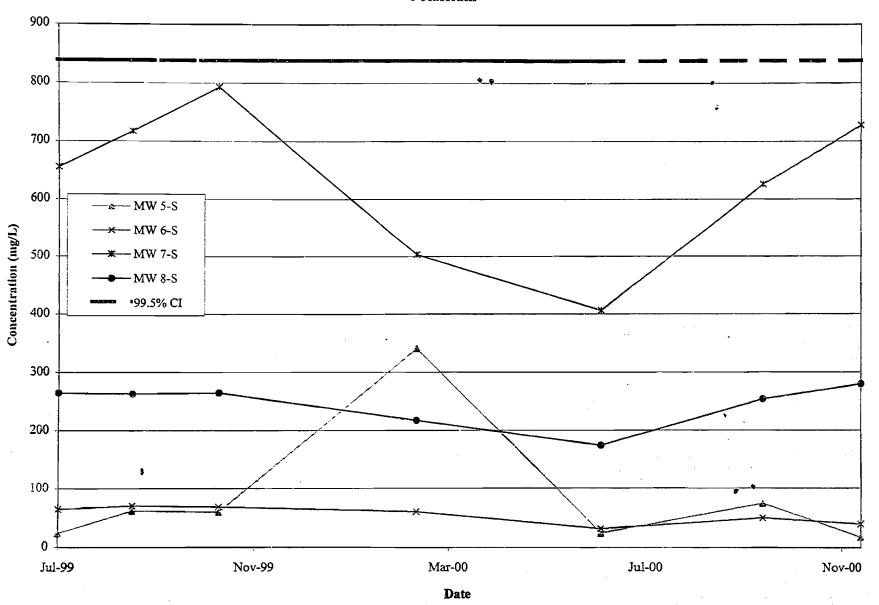


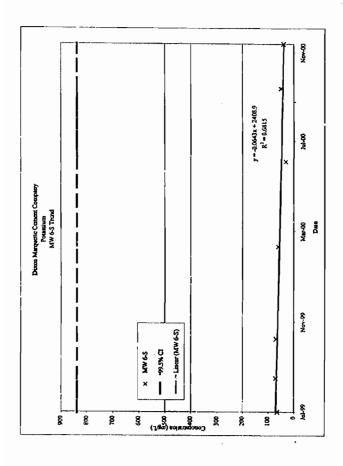


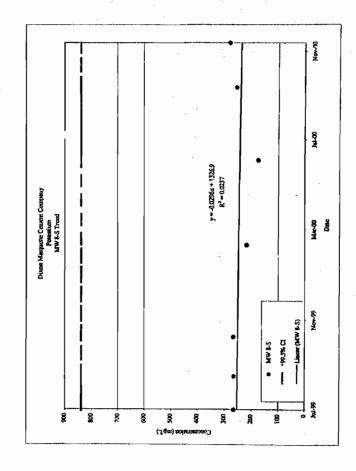


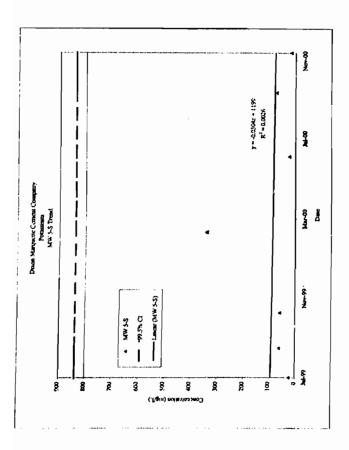


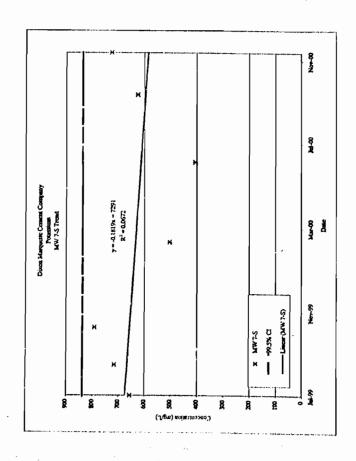
### Dixon Marquette Cement Company Potassium



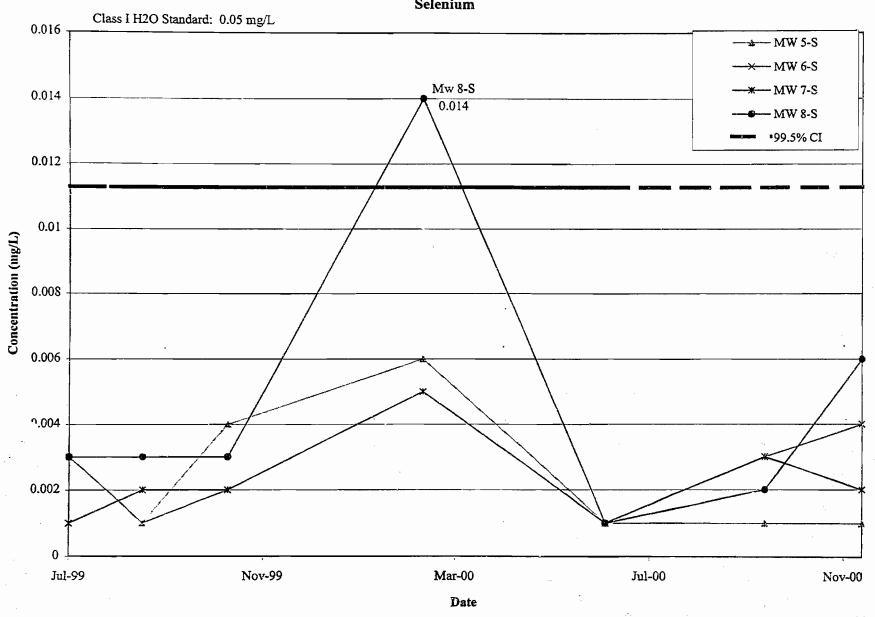


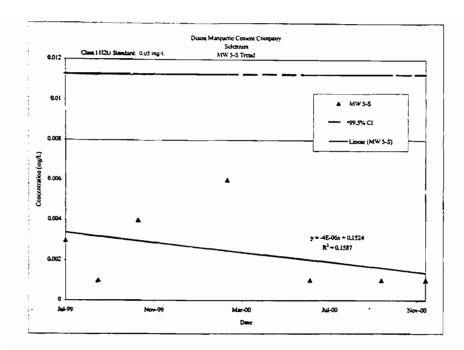


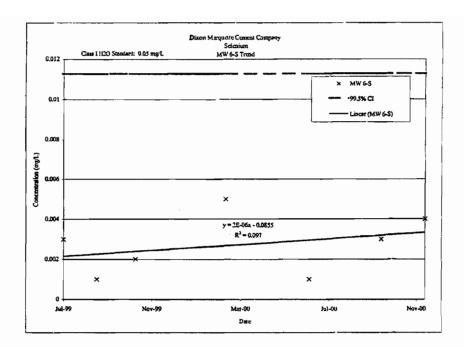


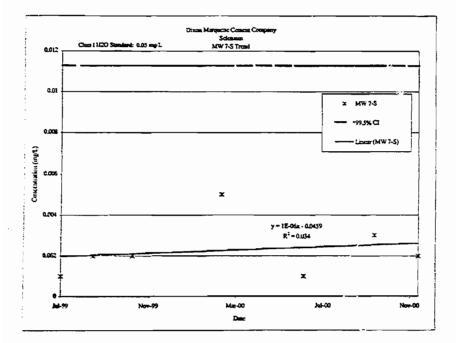


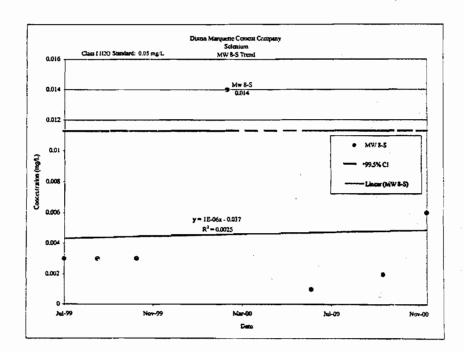
#### Dixon Marquette Cement Company Selenium



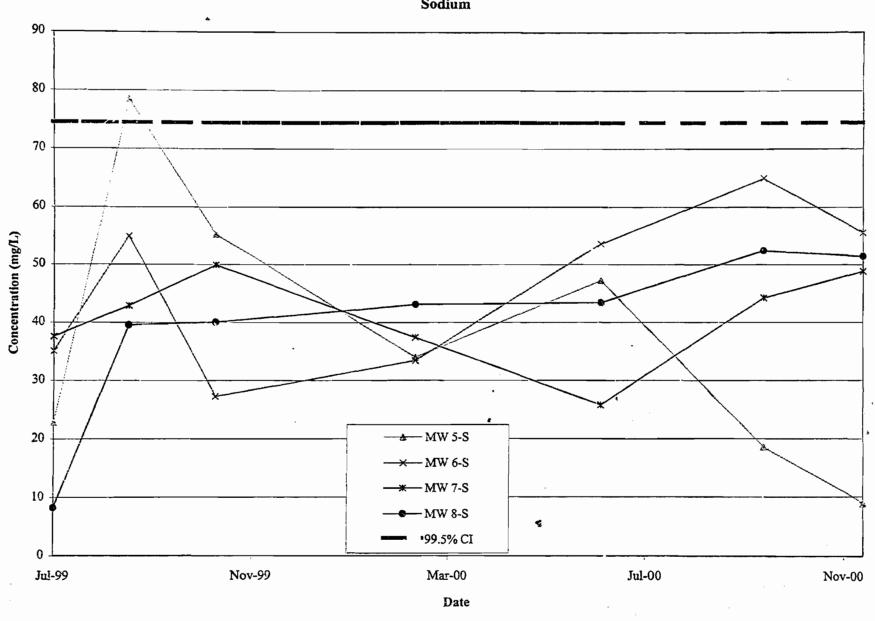


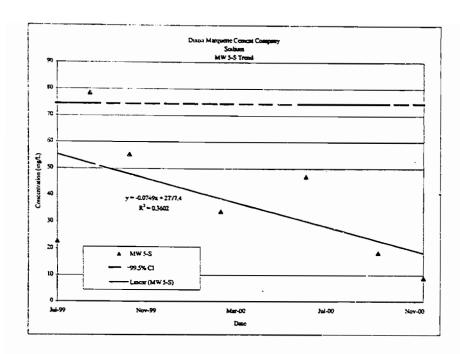


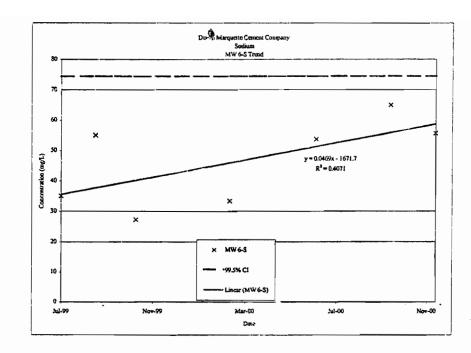


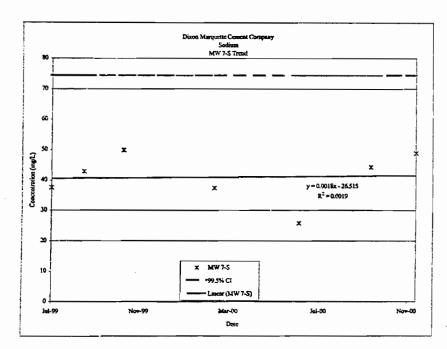


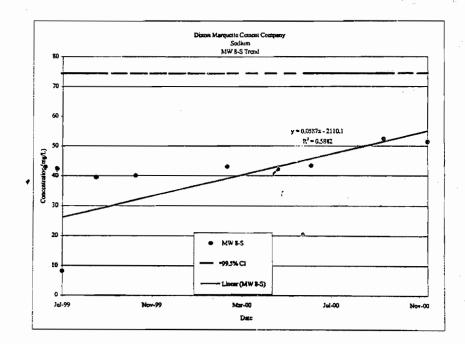
### Dixon Marquette Cement Company Sodium



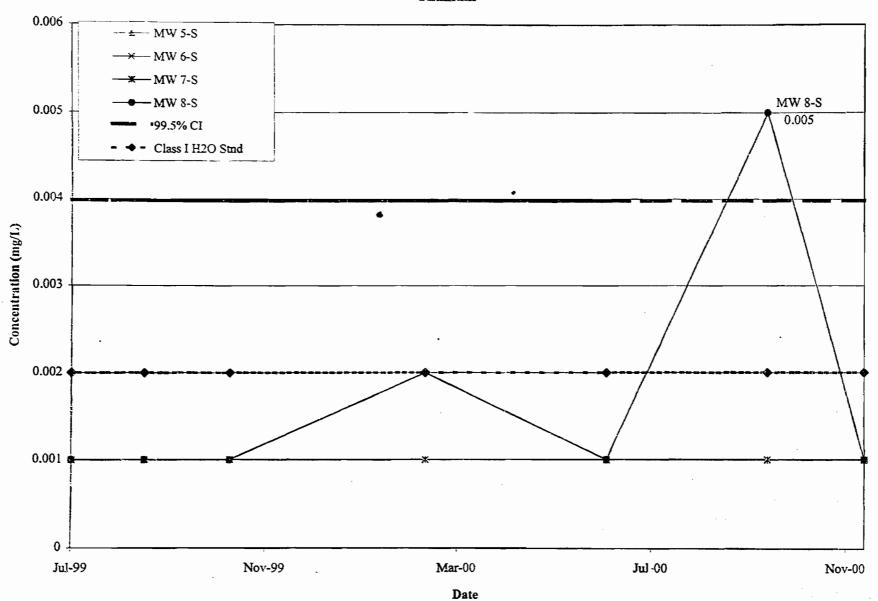


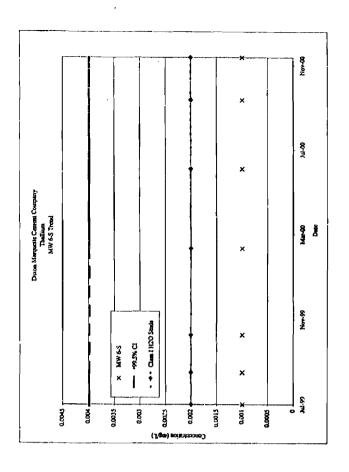


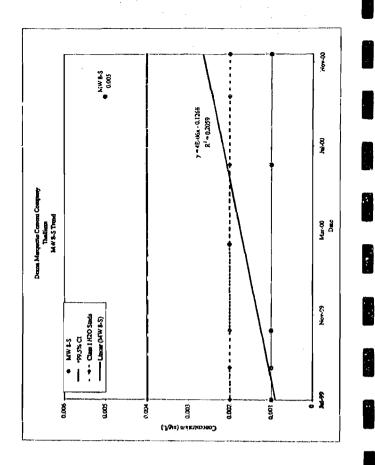




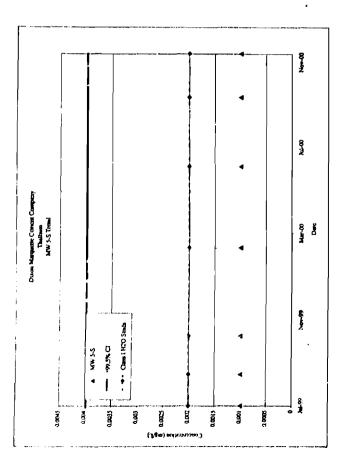
#### Dixon Marquette Cement Company Thallium

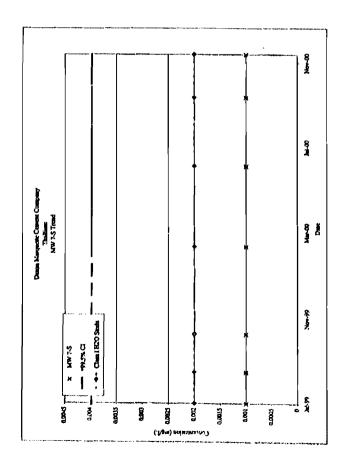






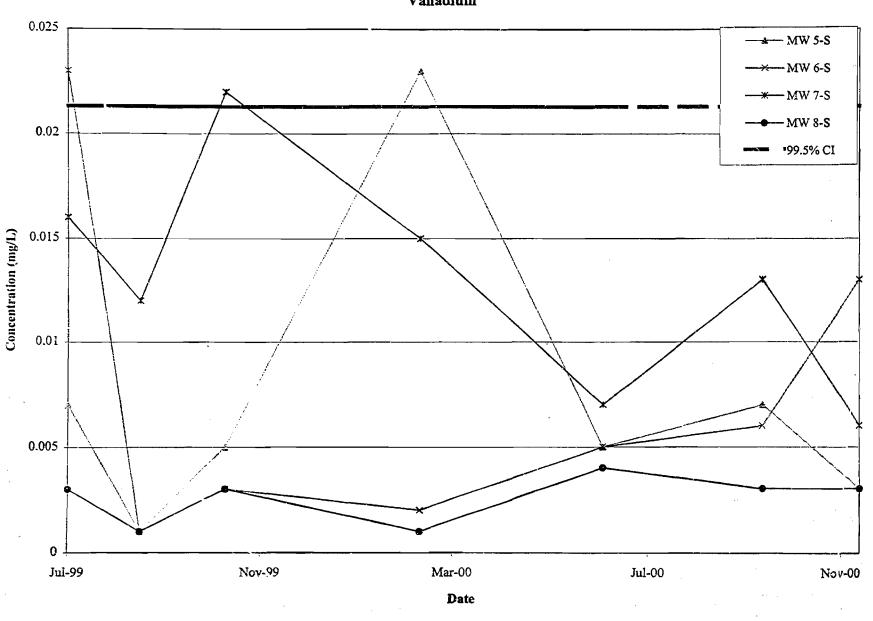
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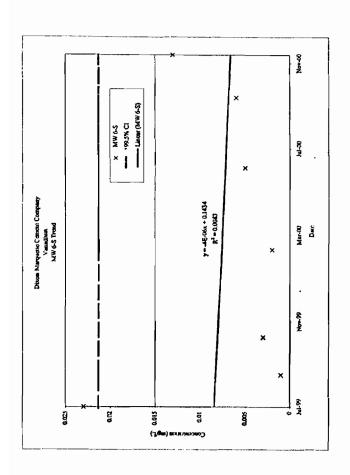


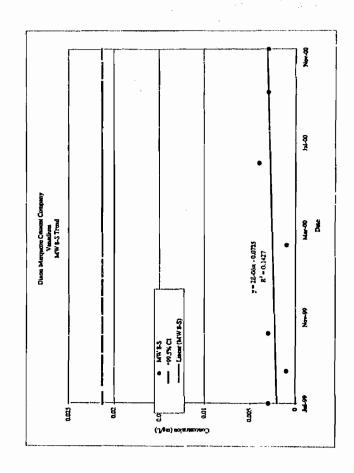


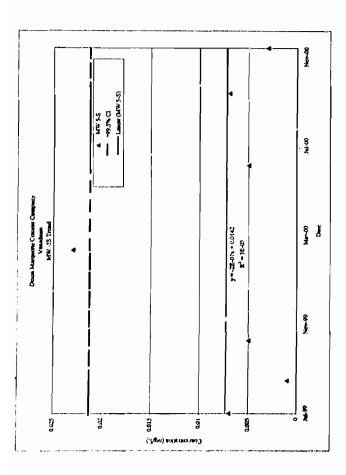
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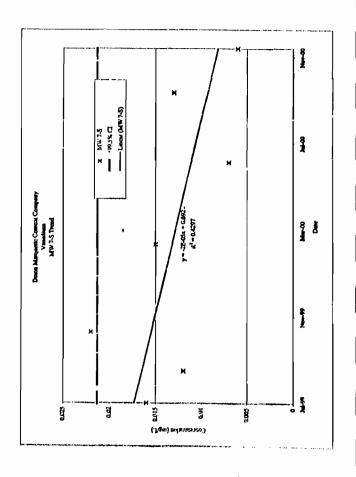
# Dixon Marquette Cement Company Vanadium



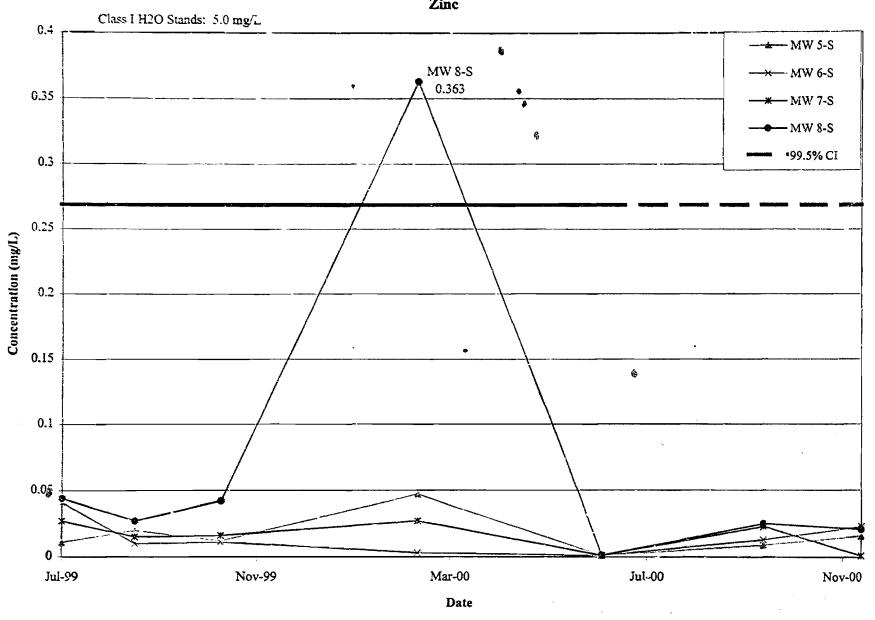


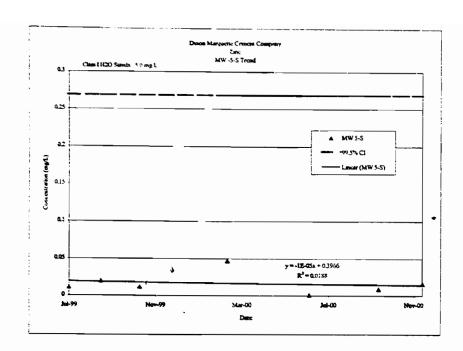


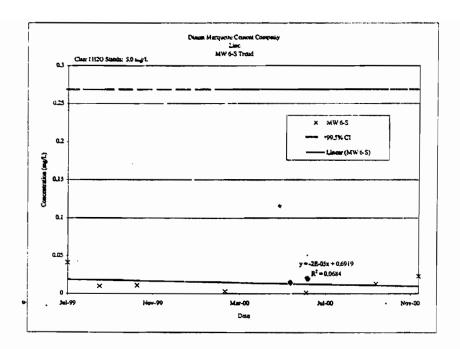


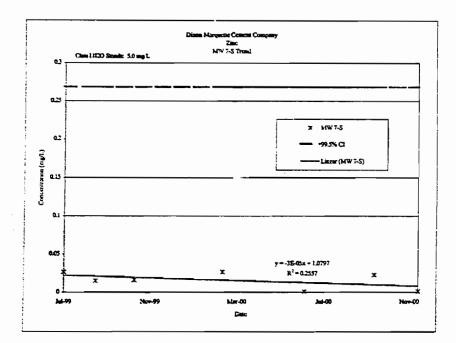


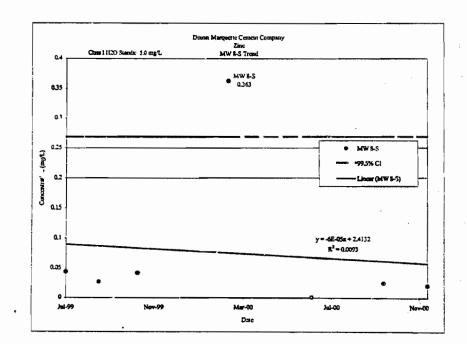
# Dixon Marquette Cement Company Zinc



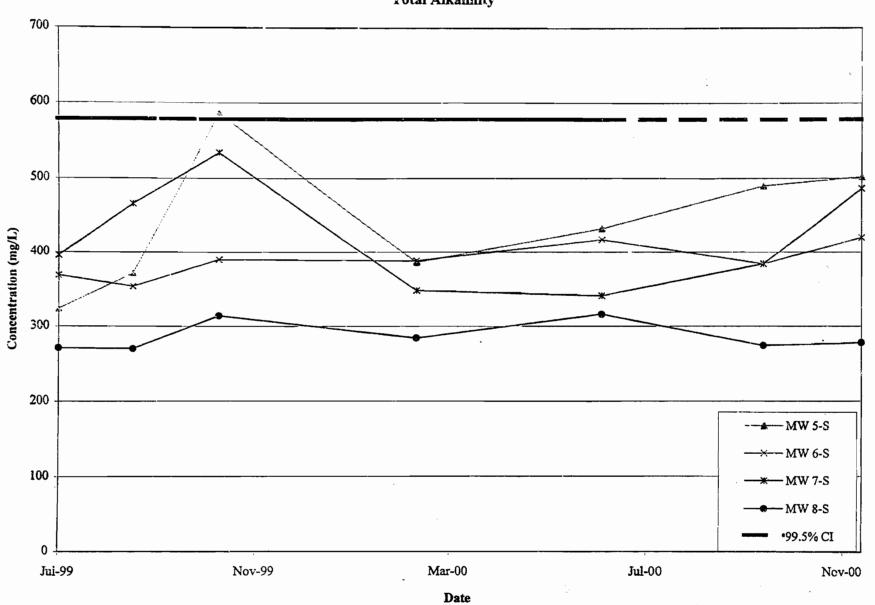


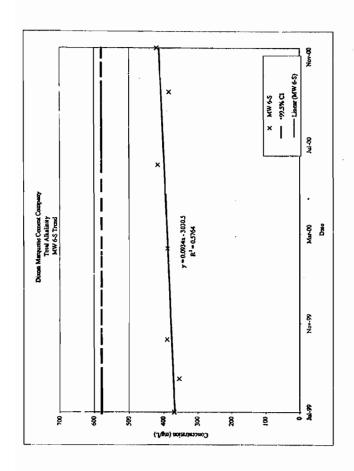


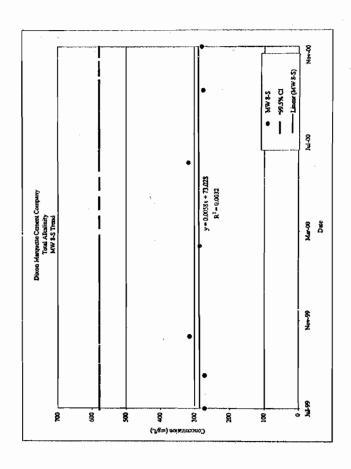


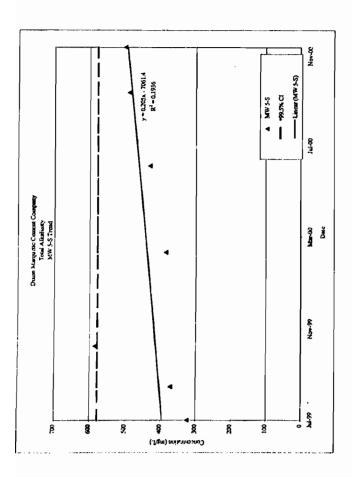


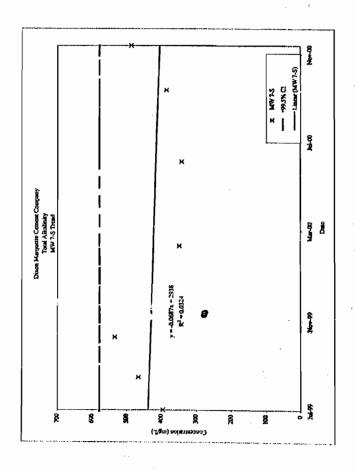
# Dixon Marquette Cement Company Total Alkalinity



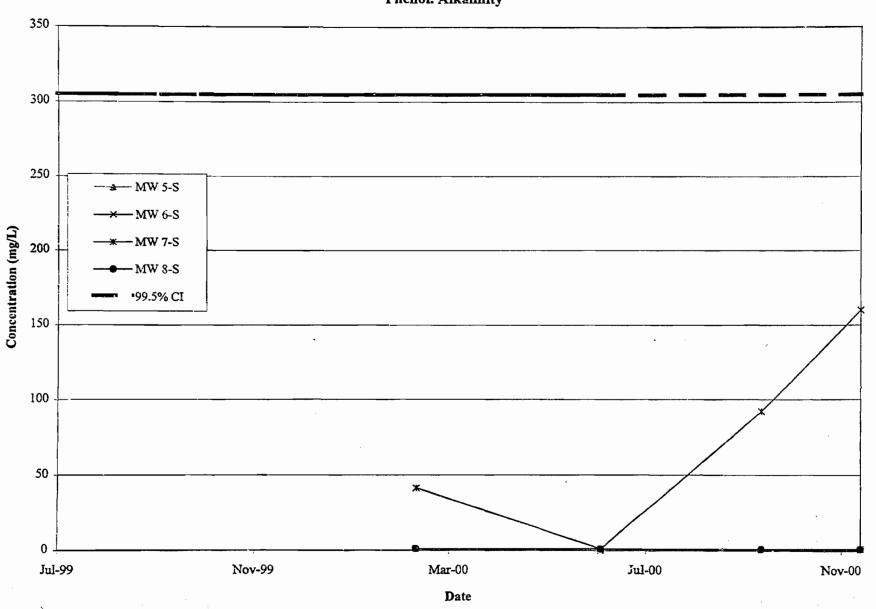


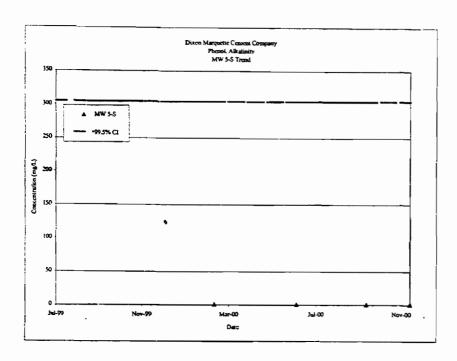


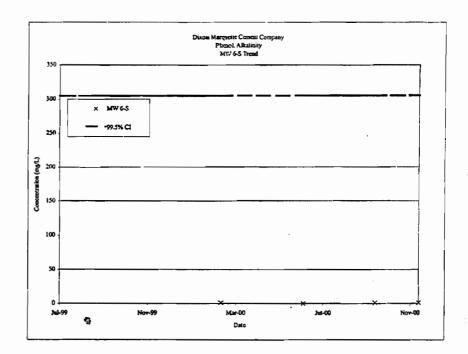


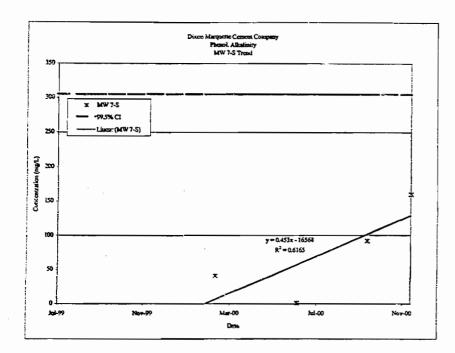


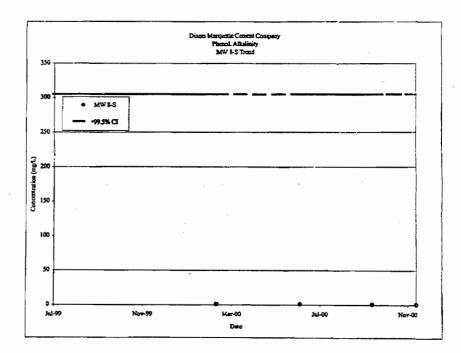
# Dixon Marquette Cement Company Phenol. Alkalinity



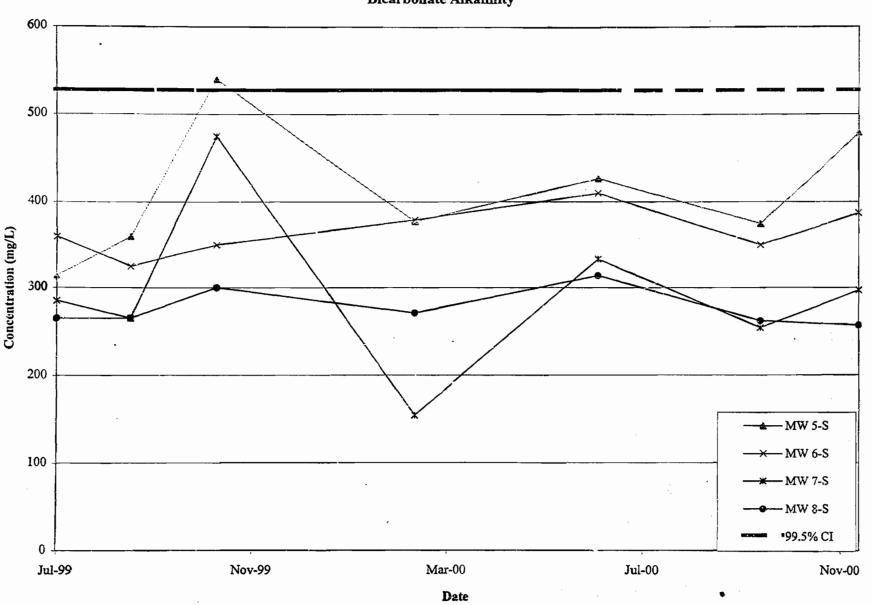


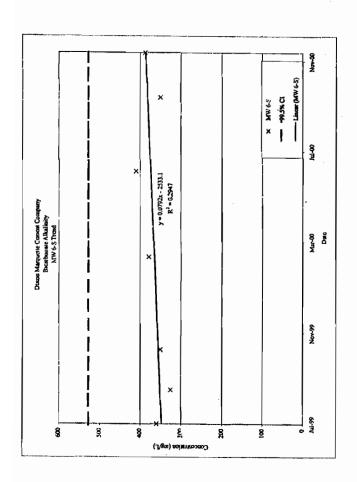


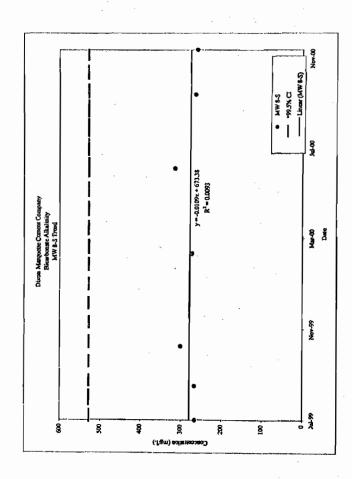


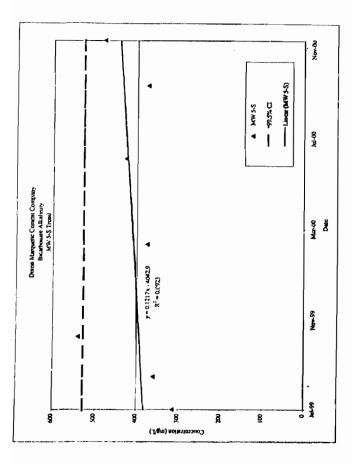


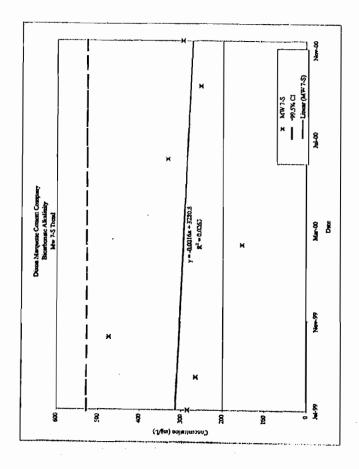
# Dixon Marquette Cement Company Bicarbonate Alkalinity



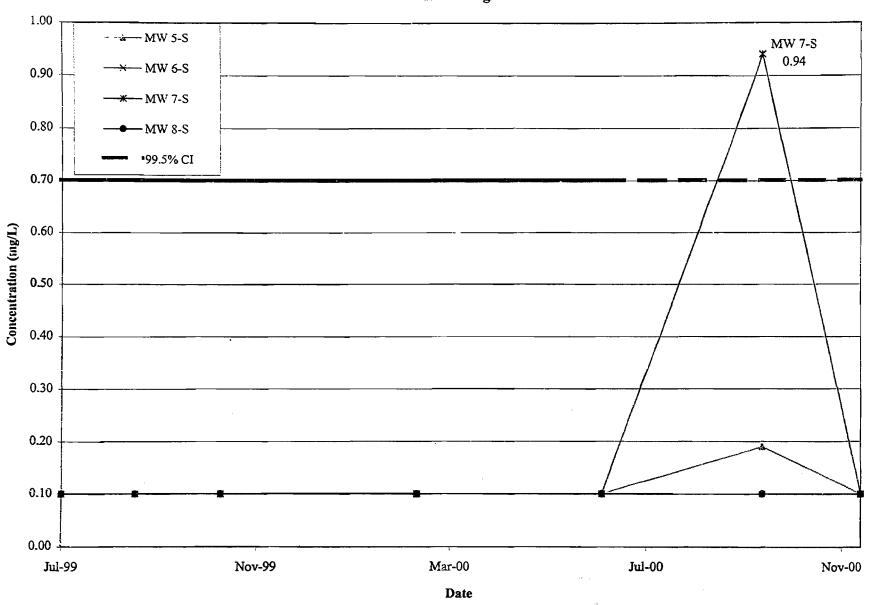


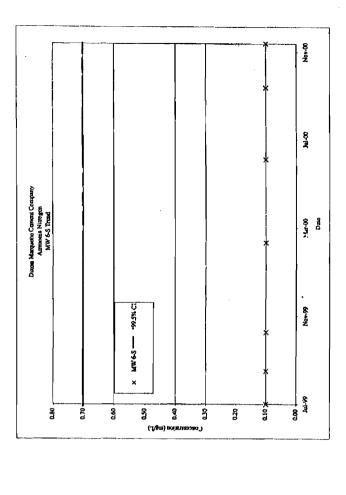


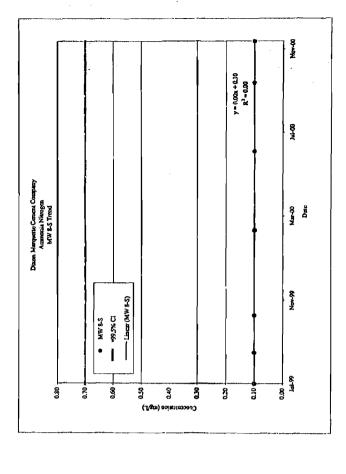


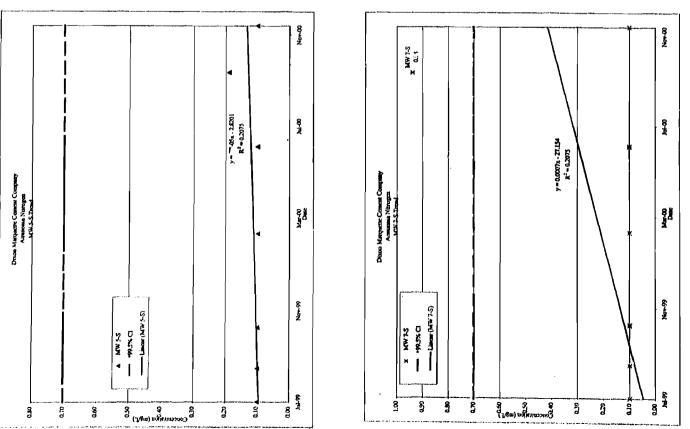


# Dixon Marquette Cement Company Ammonia Nitrogen

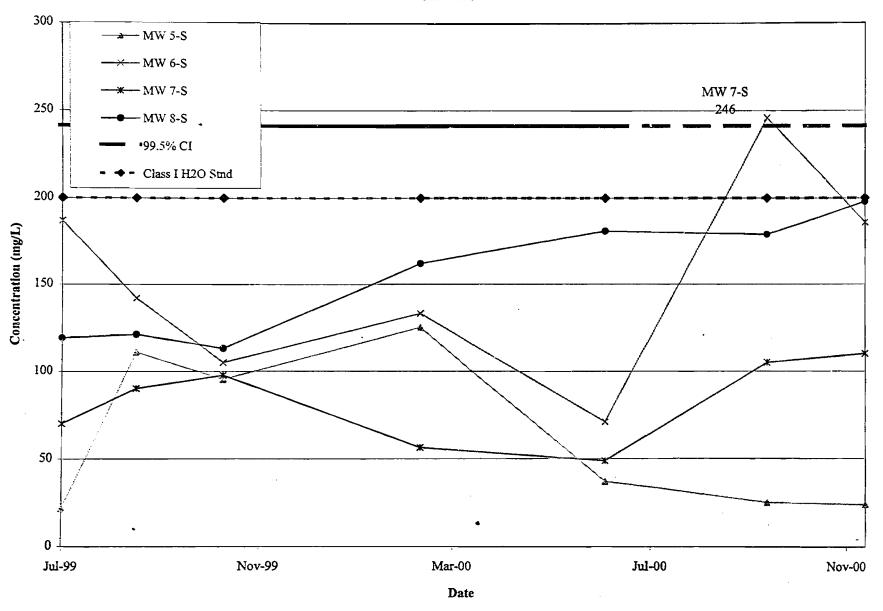


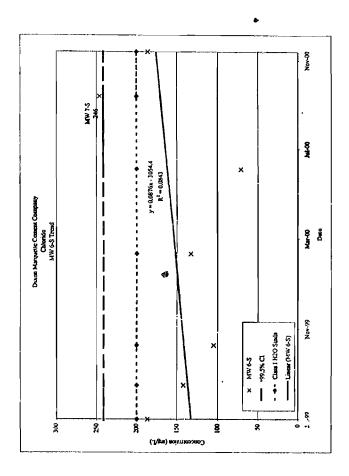


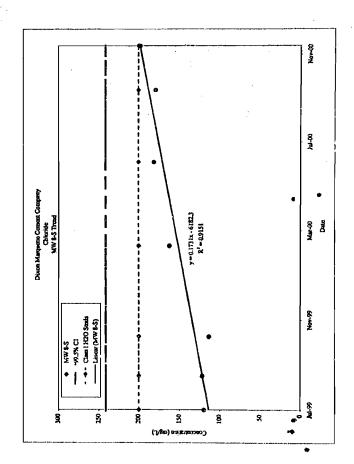


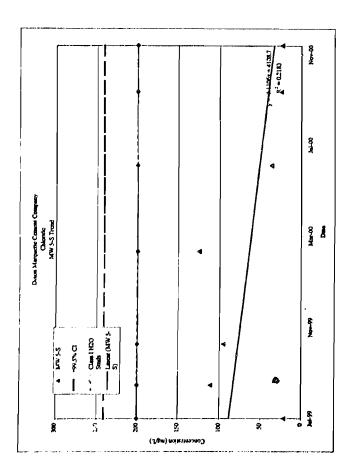


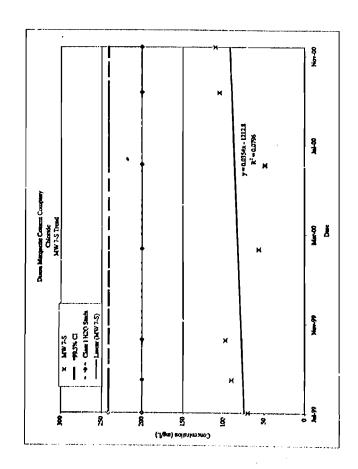
Dixon Marquette Cement Company Chloride



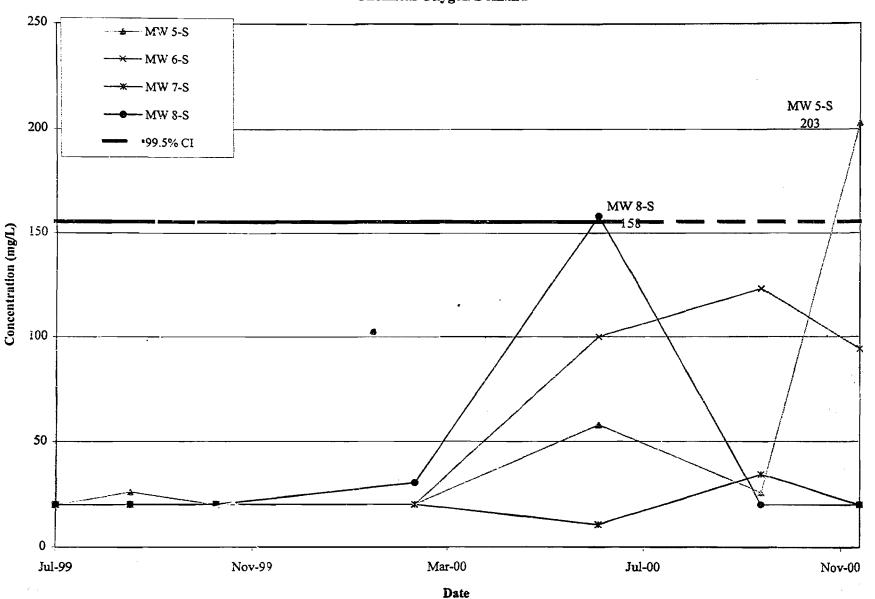


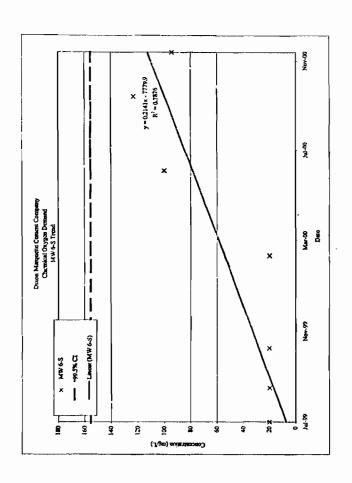


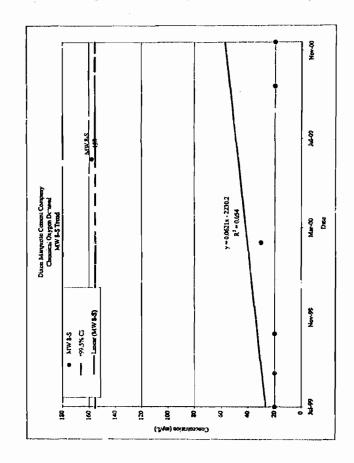


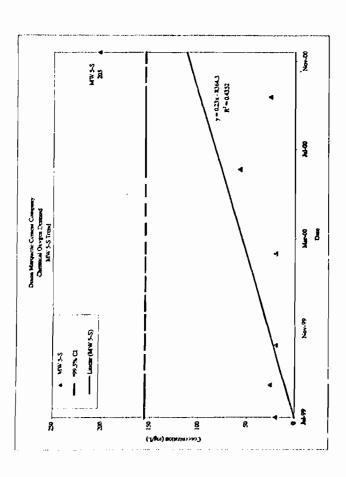


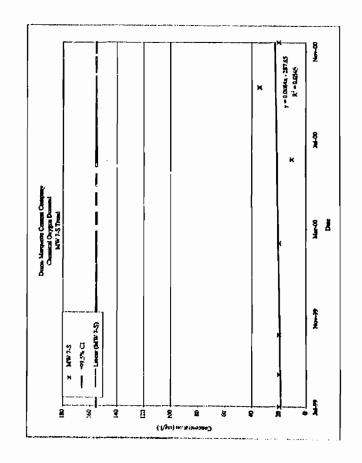
## Dixon Marquette Cement Company Chemical Oxygen Demand



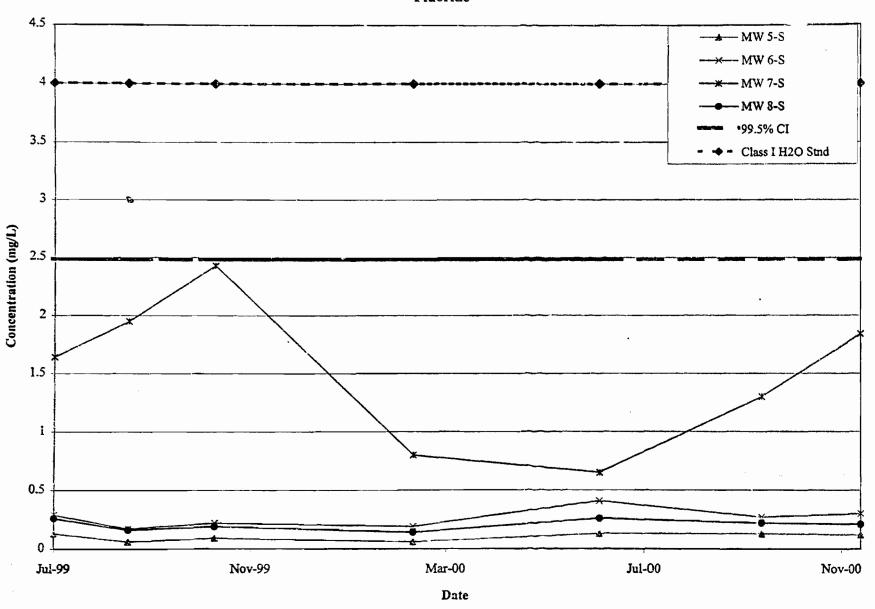


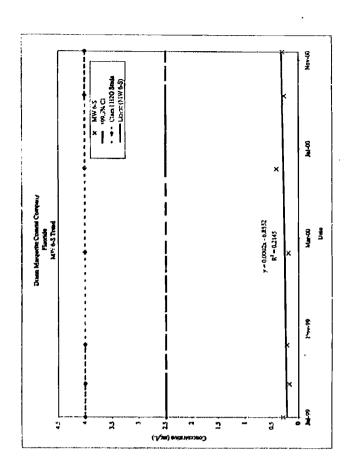


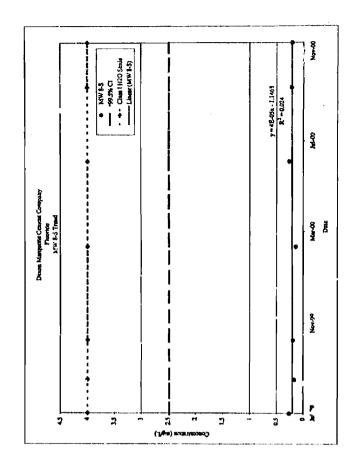




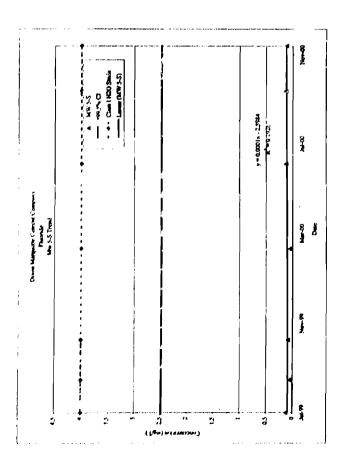
## Dixon Marquette Cement Company Fluoride

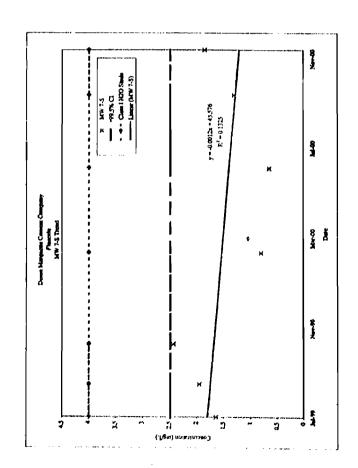




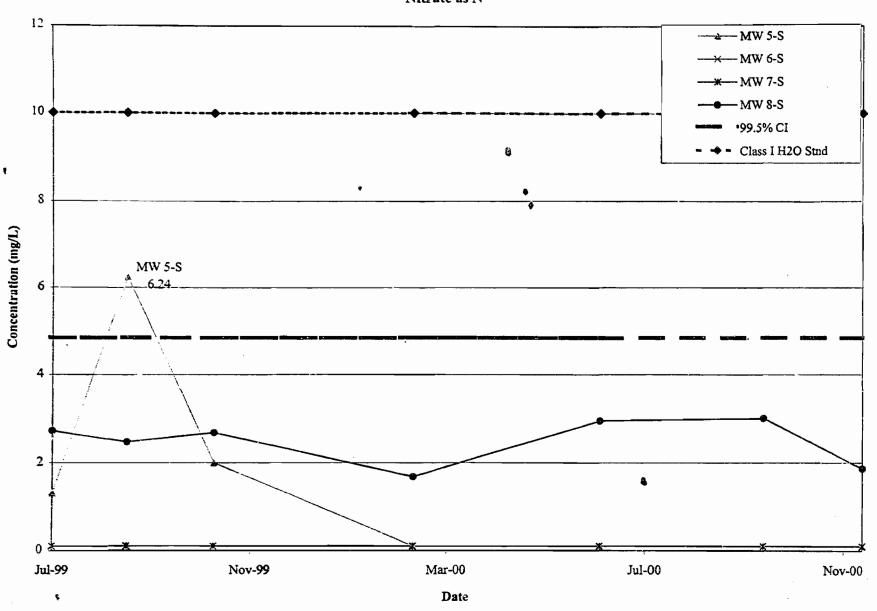


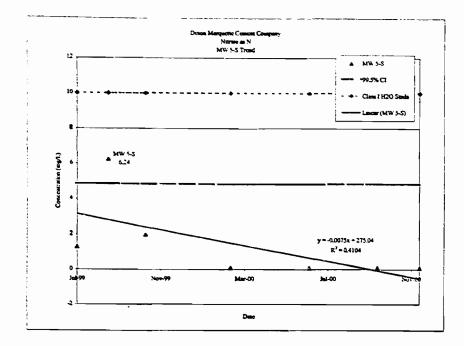
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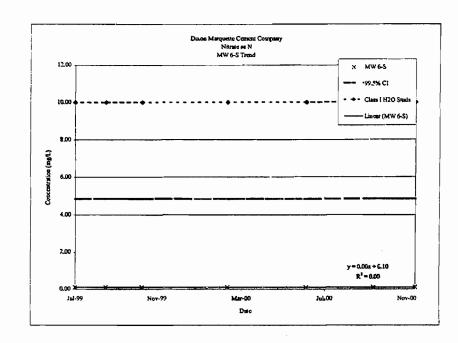


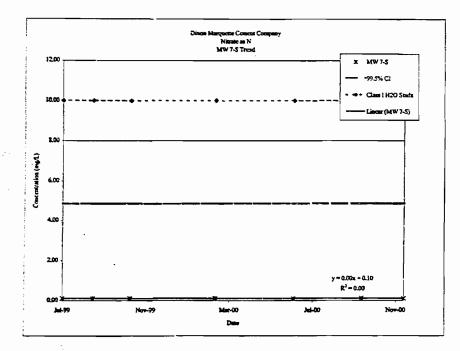


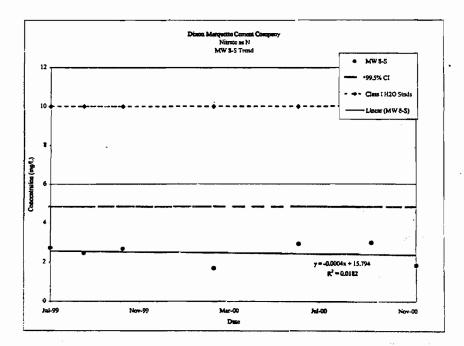
Dixon Marquette Ceme... Company Nitrate as N



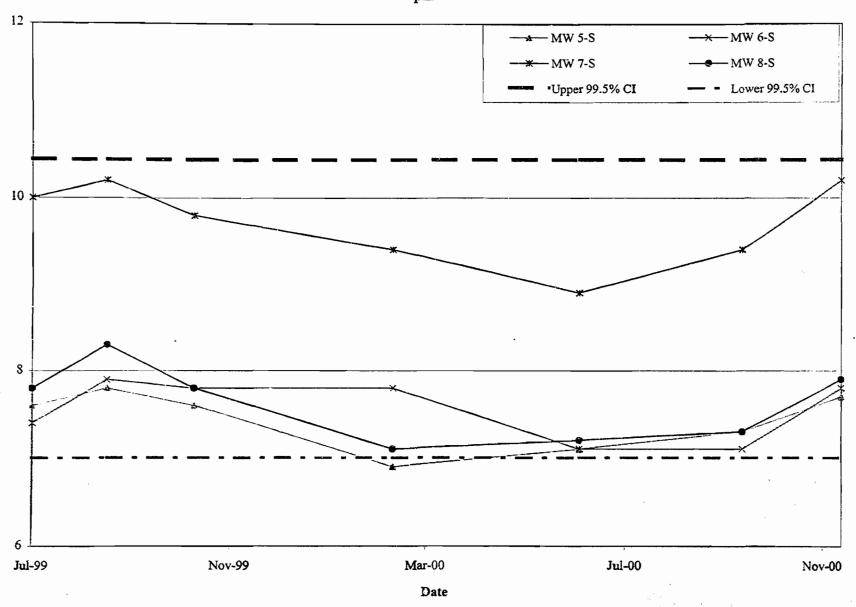


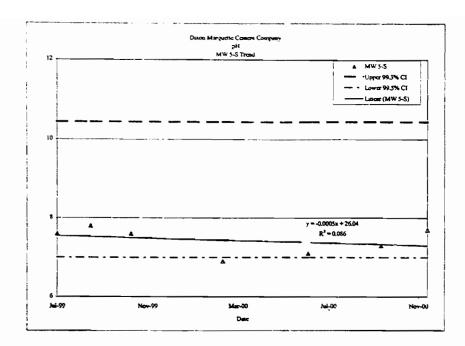


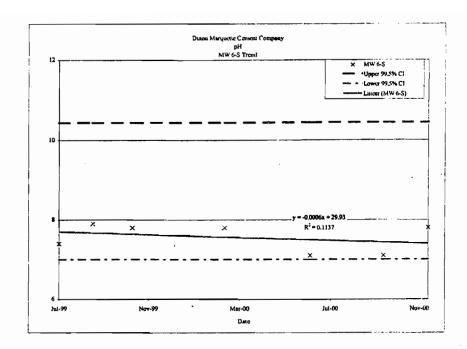


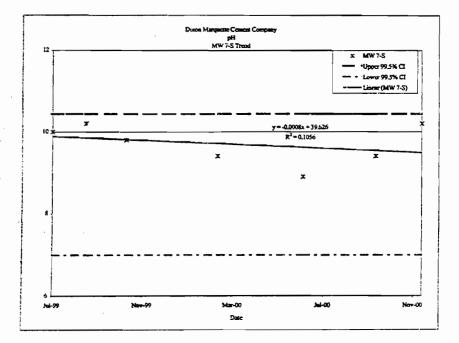


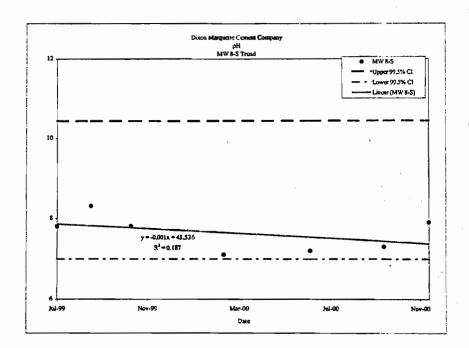
Dixon Marquette Cement Company pH



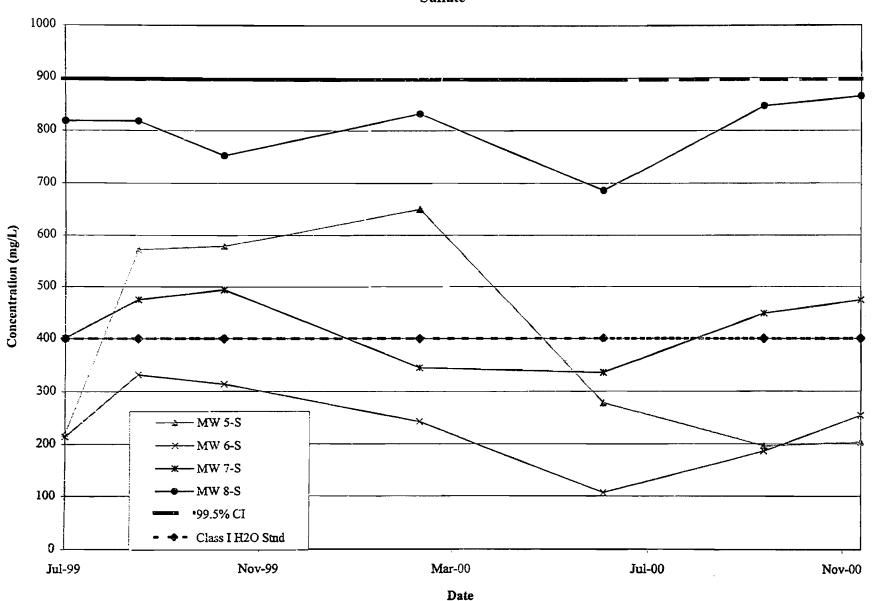


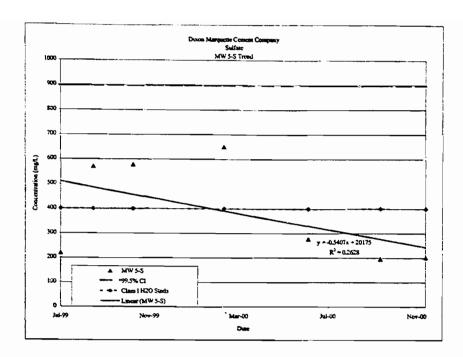


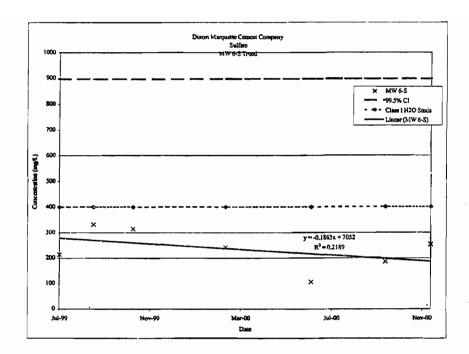


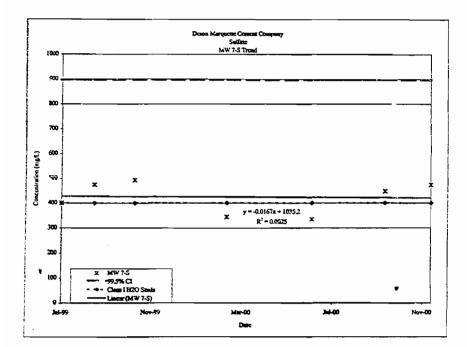


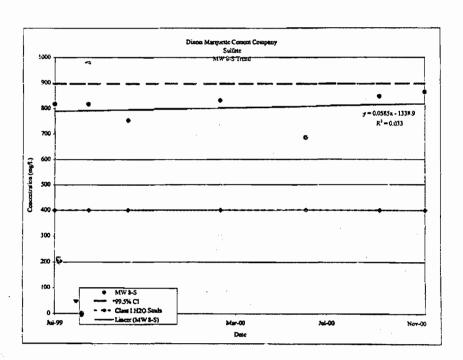
## Dixon Marquette Cement Company Sulfate



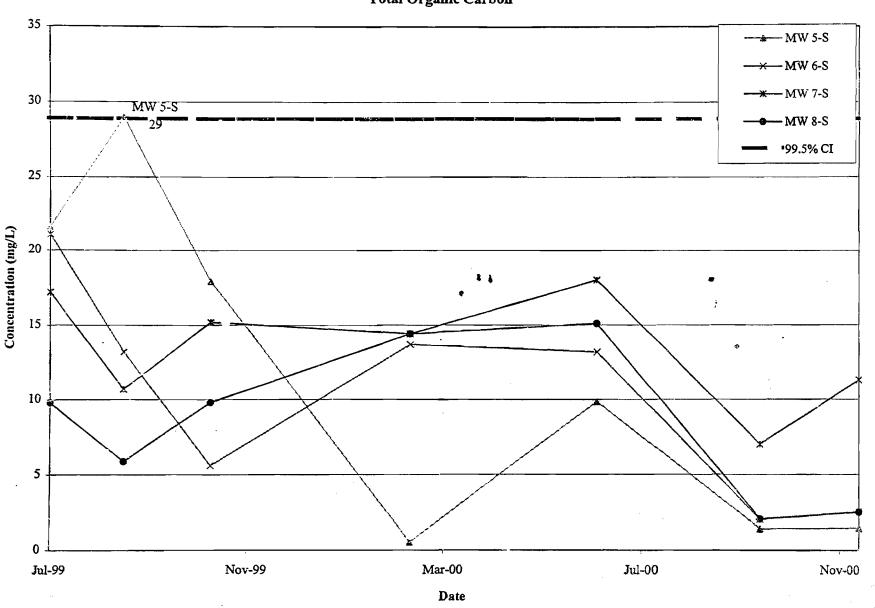


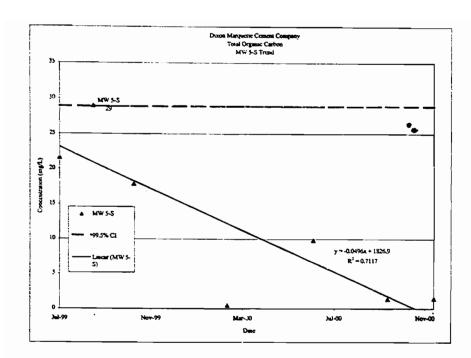


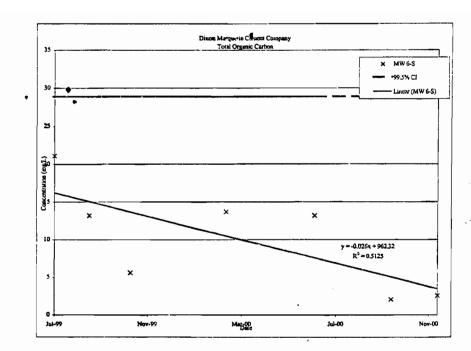


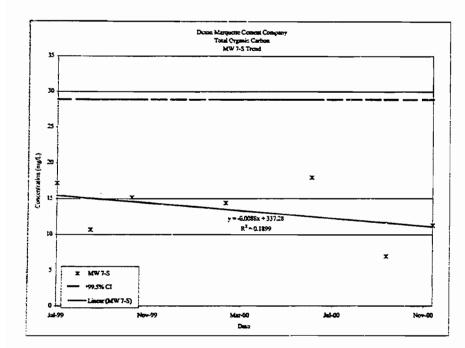


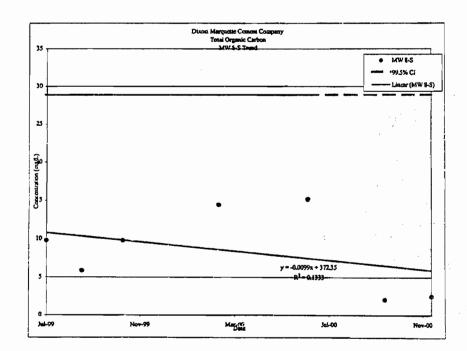
# Dixon Marquette Cement Company Total Organic Carbon



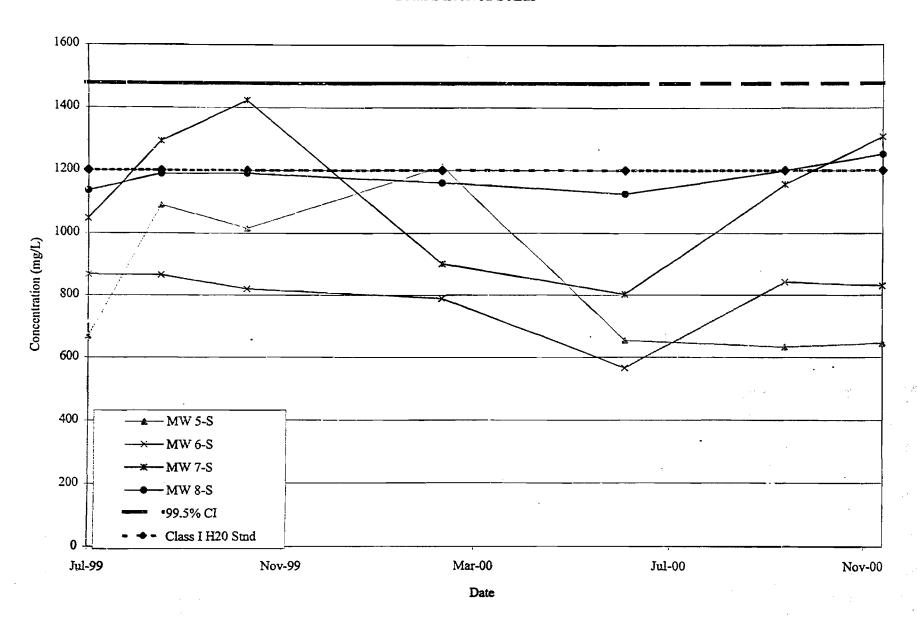


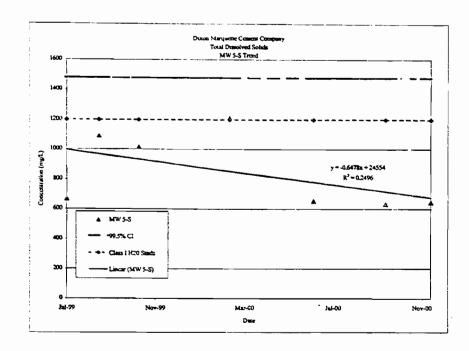


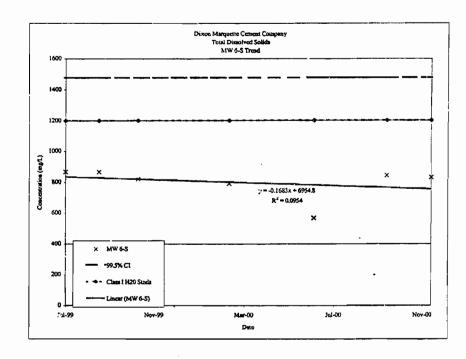


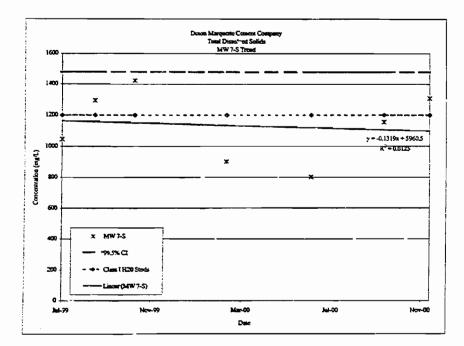


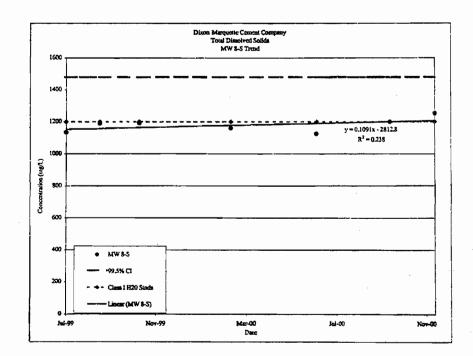
## Dixon Marquette Cement Company Total Dissolved Solids



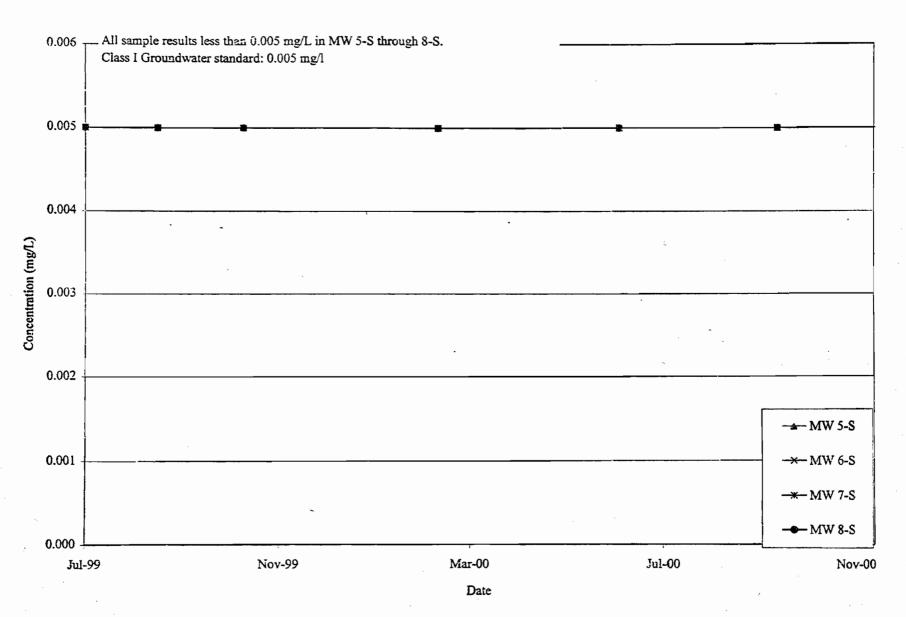




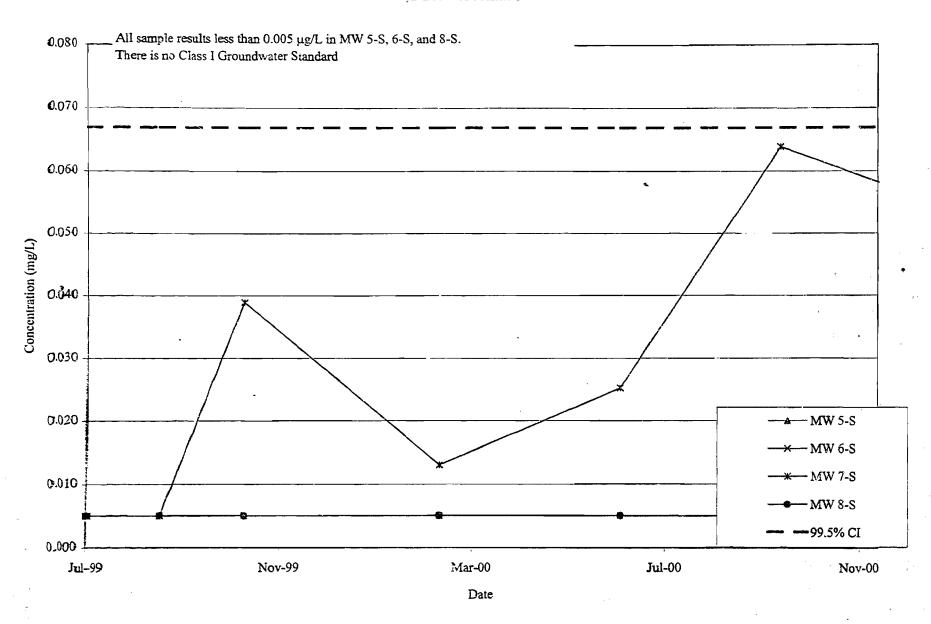


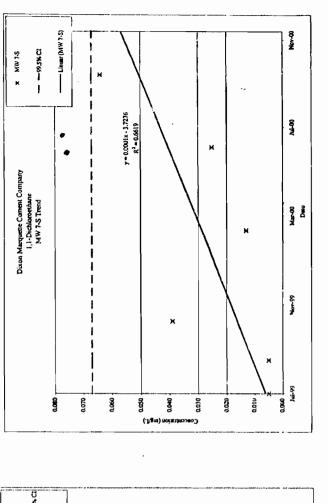


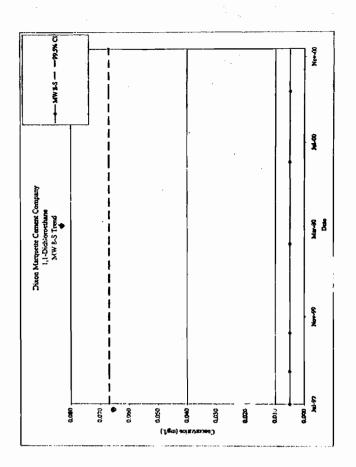
#### Dixon Marquette Cernent Company Benzene

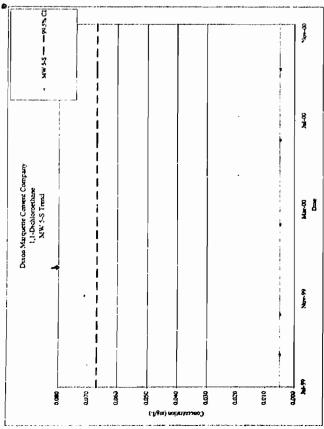


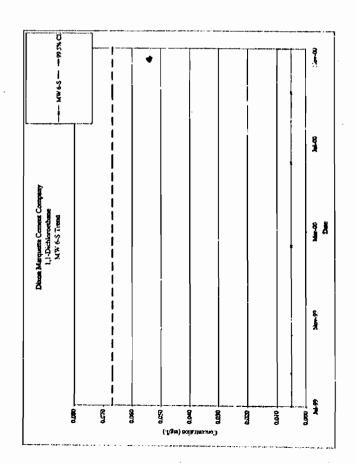
# Dixon Marquette Cement Company 1,1-Dichloroethane



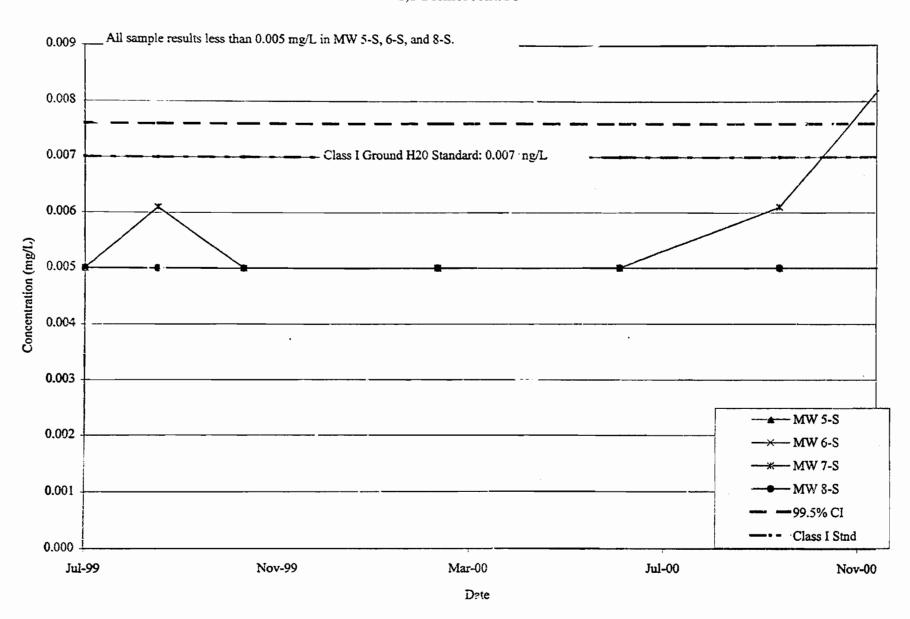


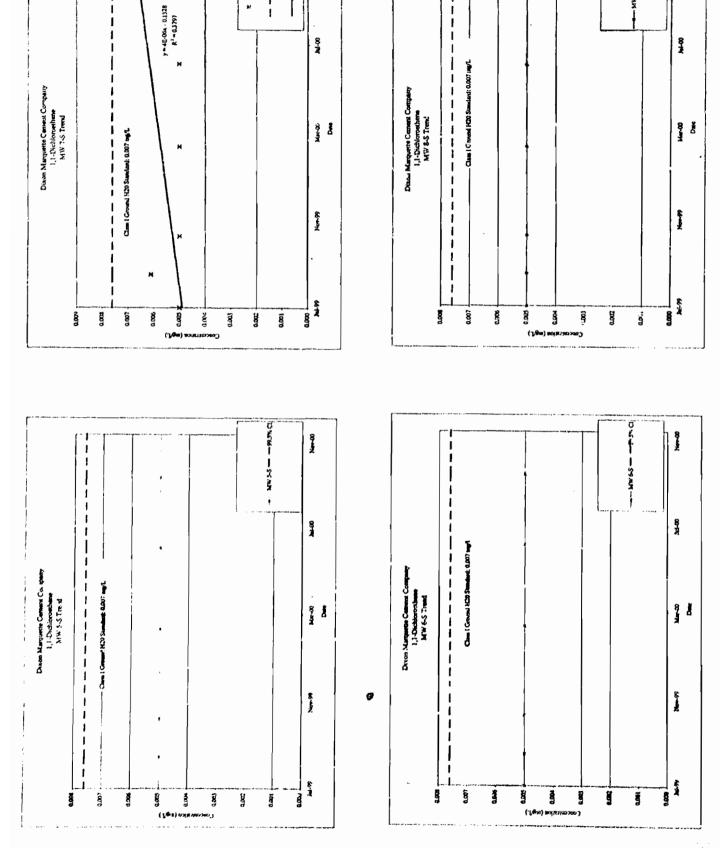






## Dixon Marquette Cement Company 1,1-Dichloroethene



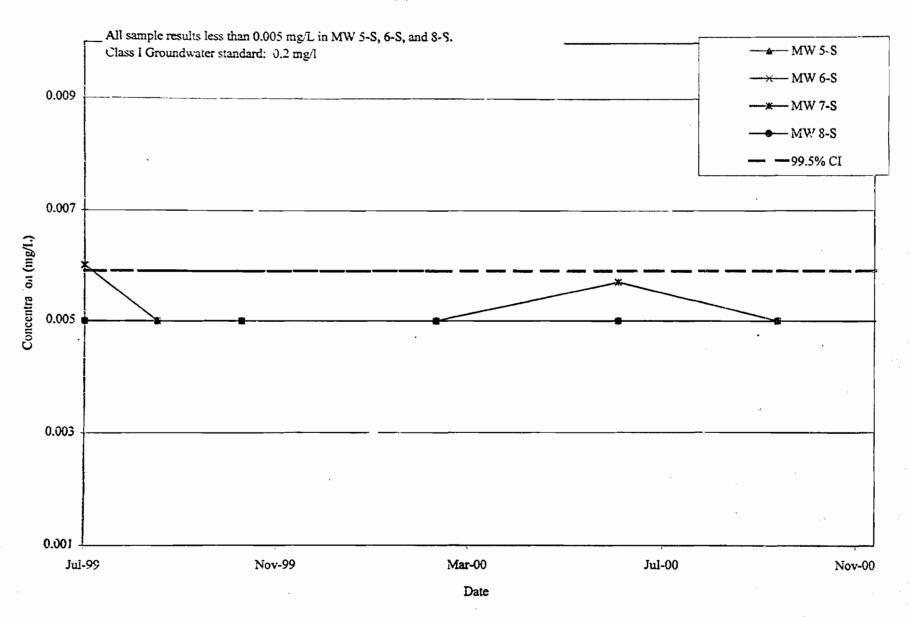


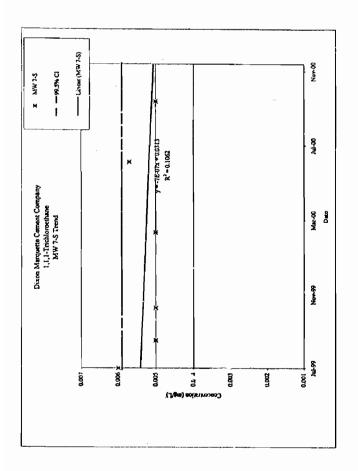
--- Linear (MW 7-S) -- -- 99.5% CI E MW7.5

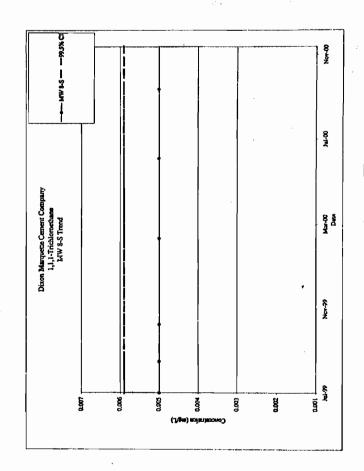
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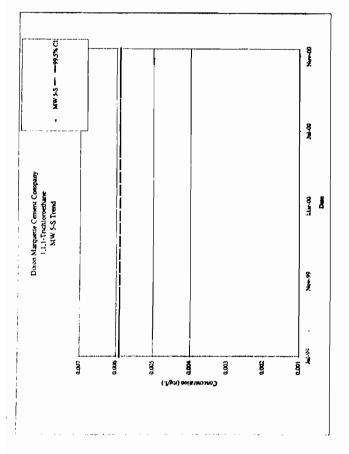
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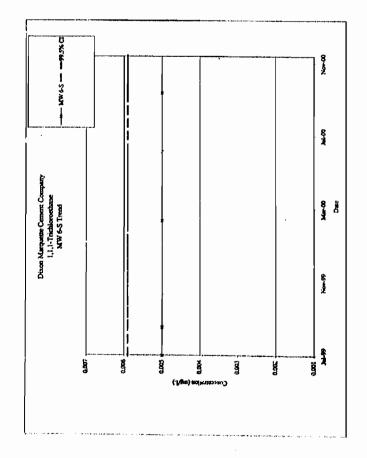
# Dixon Marquette Cement Company 1,1,1-Trichloroethane











#### Dixon Marquette Cement Company Toluene

