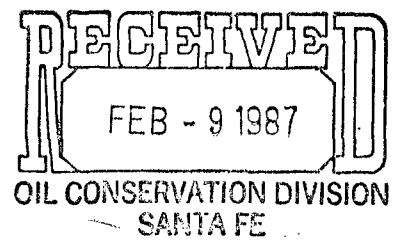


GW - 1

REPORTS

YEAR(S):

1987



A FINAL REPORT ON
SECTION 3013 ADMINISTRATIVE ORDER
WORK ELEMENTS

Prepared for
BLOOMFIELD REFINING COMPANY
Bloomfield, New Mexico

Prepared by
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CHAPTER 1

INTRODUCTION

This report submits the findings of extensive monitoring, testing, and analyses performed at the Bloomfield Refining Company, Inc. (BRC) refinery located in Bloomfield, New Mexico. Work elements described herein were established in an EPA-approved work plan to meet the directives of an administrative order issued to Gary Energy Corporation and BRC pursuant to Section 3013 of the Resource Conservation and Recovery Act (RCRA), 42 U.S.C. §6934 (Docket No. RCRA-3013-00-185).

The work plan addressing the directives of the administrative order was prepared by Engineering-Science, Inc. (ES) and submitted to EPA in July 1985. The approved work plan identifies seven separate work tasks: an electrical resistivity (ER) survey, installation of groundwater monitoring wells, groundwater monitoring well sampling, groundwater monitoring well water level measurements, groundwater monitoring well slug tests, surface water sampling, and final report. The work plan also provides an explicit protocol for execution of each of the tasks.

All work tasks with the exception of the surface water sampling task have been completed. BRC staff have performed the Hammond Ditch sampling and the monitoring well sampling and level measurements, and have administered to the installation of monitoring wells 7 through 10. ES has performed the electrical resistivity survey and monitoring well slug tests. These tasks were performed consistent with the specifications in the approved work plan. A time extension has been requested for the surface water sampling task to permit sampling of the San Juan River during representative low-flow conditions (heretofore nonexistent during the 13-month work period).

In summary, BRC has executed each of the work elements of the 3013 administrative order in a timely manner and consistent with guidelines in the approved work plan. This report, which fulfills the requirements of the administrative order, presents findings of these work elements which

cumulatively suggest there are generally low contaminant levels found at the refinery. In the information that follows, chapters 2 through 7 address each of the specific work elements as required by the approved work plan. Chapters 8 and 9 contain conclusions and recommendations based on these findings.

An amendment to this report will be submitted with results of the San Juan River sampling effort, which are outstanding due to continued high flow rates. Upon submission of the amendment, the requirements of the 3013 administrative order will have been completed.

CHAPTER 2

ELECTRICAL RESISTIVITY SURVEY

BACKGROUND

The purpose of the electrical resistivity (ER) survey at the Bloomfield refinery was to develop a data base that would assist in determining the location, areal extent, and direction of migration of subsurface hydrocarbons. The survey, which was conducted during the period January 27 to February 7, 1986, in a manner consistent with the EPA-approved work plan, accomplished this purpose. Findings of the survey were sent to Mr. Steven Schwartz of EPA Region VI in February 1986 in a formal report entitled "Electrical Resistivity Survey," Bloomfield Refining Company.

METHODOLOGY

Since the ER survey technique is an indirect measurement method that reports relative changes in the electrical signature of subsurface features, the process of obtaining measurements must therefore refer to known subsurface features. Interpretations of these data are based heavily on correlation of ER measurements with verified locations and depths of known subsurface features such as subcrops, seeps, water, and, in this case, subsurface hydrocarbons. The success of the ER survey is therefore contingent on the degree to which the ER measurements can be verified to correlate with known subsurface features.

During the survey, both vertical (sounding) and horizontal (profile) measurements were made to develop the data base. These measurements were made using a Bison Model 2530B earth resistivity meter together with published techniques for performing the measurements referenced in the ER report. Investigators utilized monitoring well logs that depicted stratigraphy with respect to depth; subcrop locations of stratigraphic layers witnessed from a bluff; the location of seeps and ponds of water; the locations of visible hydrocarbon seepage in Hammond

Ditch and on the bluffs adjacent to the San Juan River; and the depth of floating hydrocarbons in monitoring well 4 (free hydrocarbon phase noted by Bloomfield staff during well samplings throughout 1986) to interpret relative changes in measurements made during the survey. The locations of measurement points are presented in Figure 2.1. Computer-enhanced interpretations of the profile measurements were generated for Figures 2.2 through 2.8. These plots provide an areal interpretation of profile ER readings for selected reference depths.

Interpretation of the sounding measurements was developed in the form of plots of apparent resistivity (measured values of ER) versus inner electrode spacing. The cumulative resistivity values displayed on these plots provide a measure of the degree of consistency of subsurface features and conditions. These plots, together with a location map showing the location of each sounding measurement, are presented as Figures 2.9 through 2.15.

SITE GEOLOGY

Both the sounding and profile measurements presented in the report via subsurface cross sections showed the geology underlying the refinery property to be generally homogeneous and to consist of four distinct geologic units. These units are the Nacimiento Formation; a cobbles and pebbles unit; a sand, silt, and clay unit; and an alluvial fill unit. Two geologic cross sections were prepared using sounding data together with data from monitoring well logs to provide an understanding of the orientation of the geology through the BRC site. The sections were selected based on site constraints and the objective of obtaining an approximately perpendicular arrangement of the sections. The sites selected are shown on Figure 2.9 as locations A-A' and B-B'. These sections, presented by Figures 2.16 through 2.17, show the geology tends to conform to a northerly dip in the top of the Nacimiento Formation and that the surface of this formation undulates.

As suggested by Figure 2.17, the slope of the Nacimiento is not continuous. Figure 2.18, developed in a discharge plan submitted to

FIGURE 2.1
BLOOMFIELD REFINING CO.: PROFILE LOCATIONS

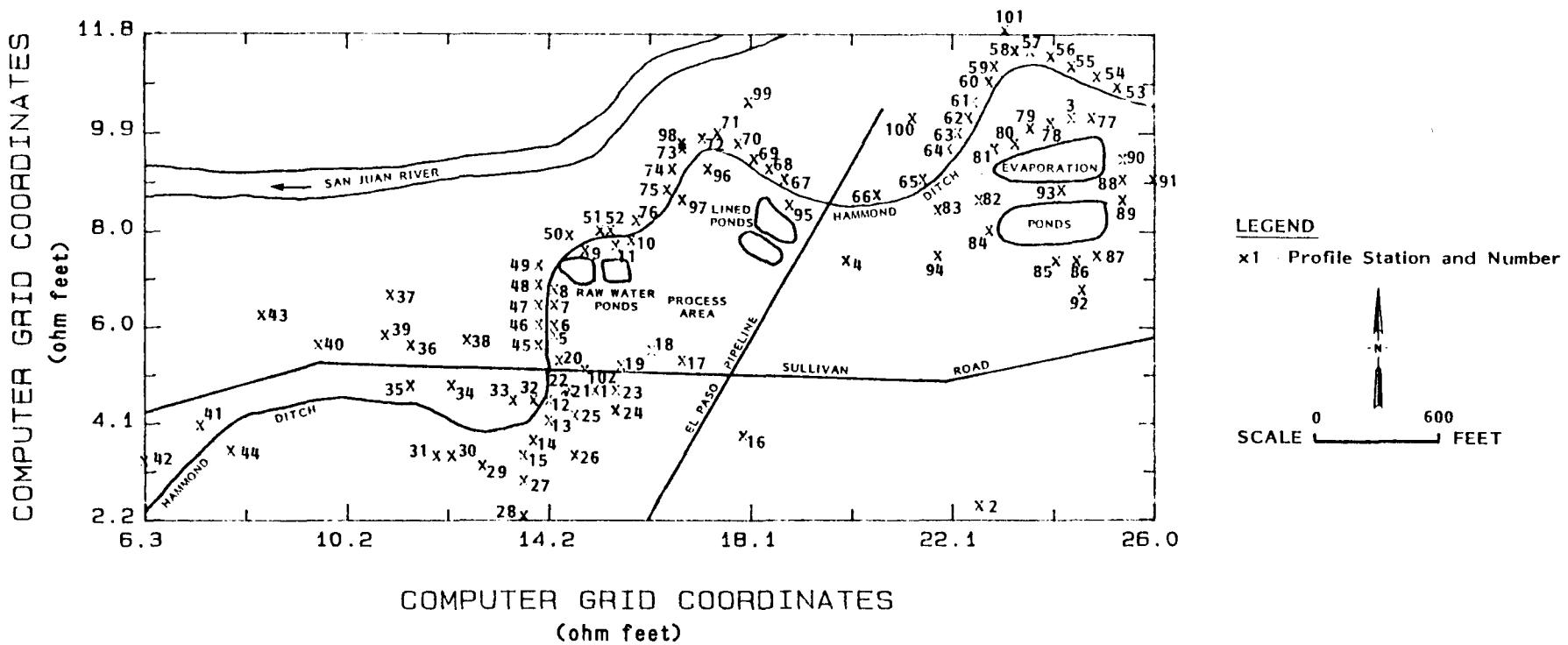


FIGURE 2.2

BLOOMFIELD REFINING CO.: 10-FT. ER PROFILE MAP

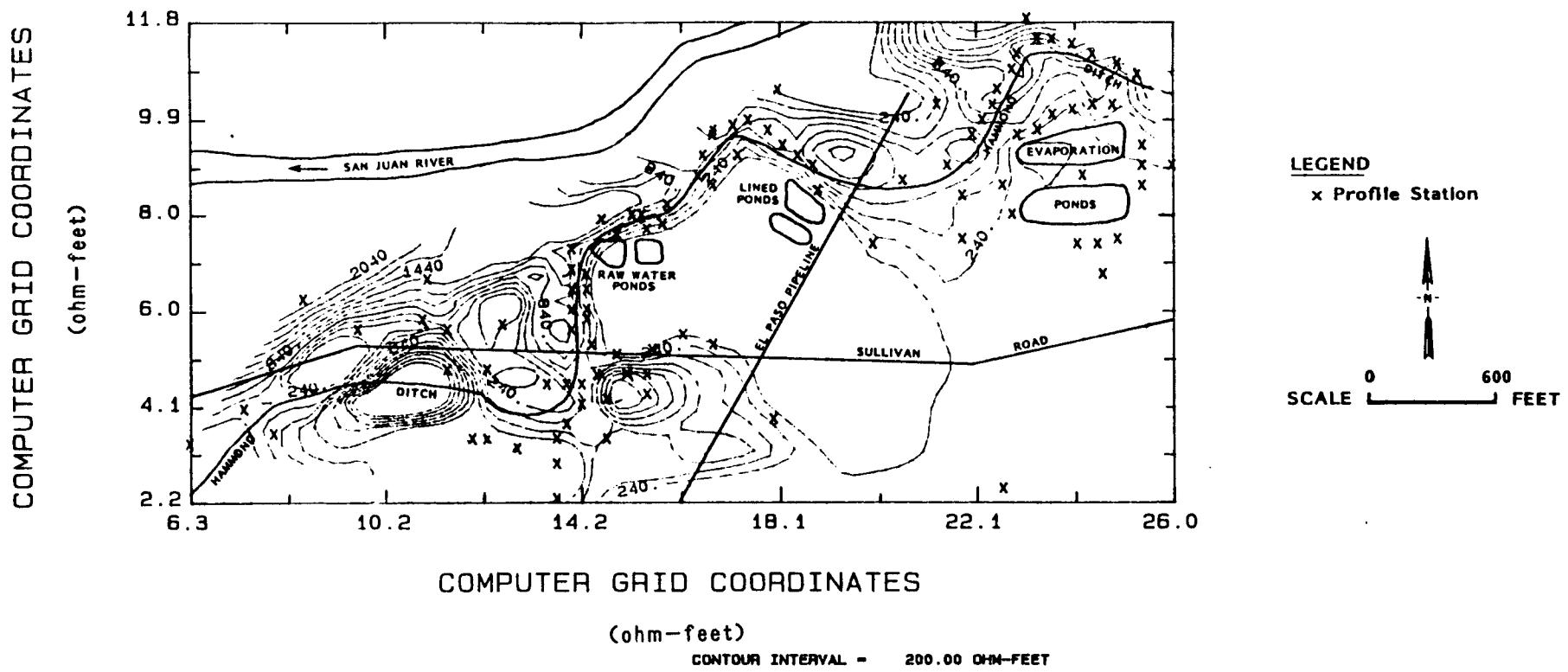
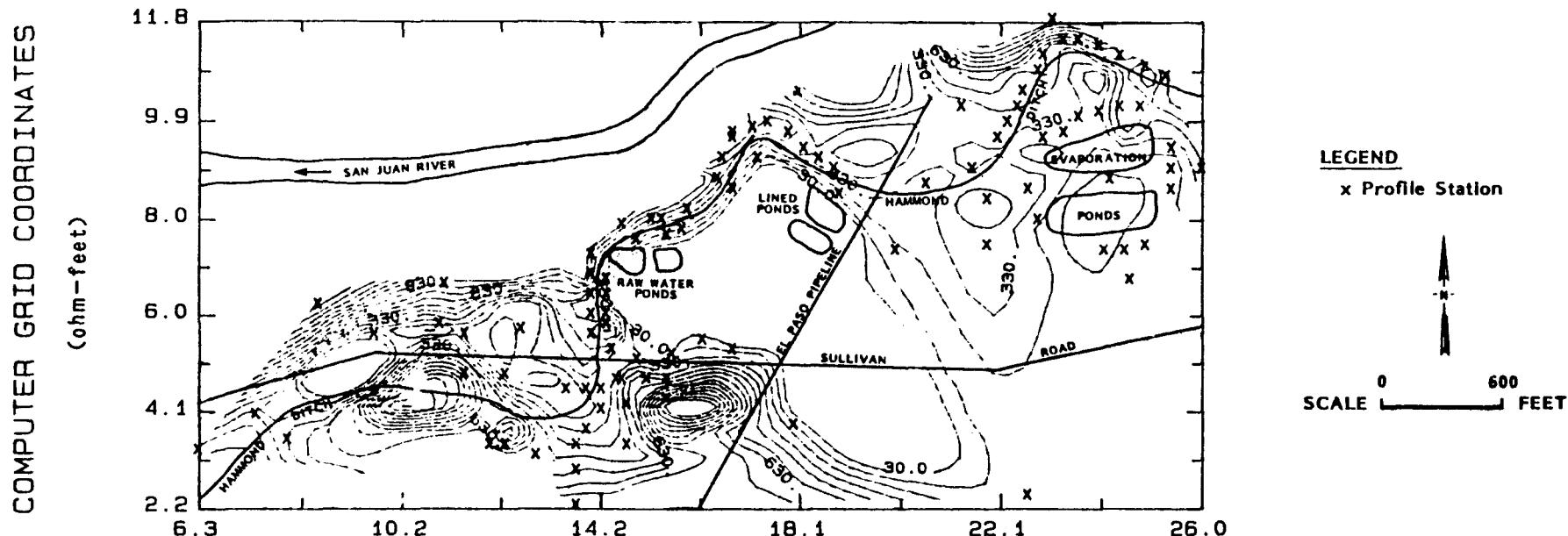


FIGURE 2.3

BLOOMFIELD REFINING CO.: 20-FT. ER PROFILE MAP



COMPUTER GRID COORDINATES

(ohm-feet)

CONTOUR INTERVAL = 100.00 OHM-FEET

FIGURE 2.4

BLOOMFIELD REFINING CO.: 30-FT. ER PROFILE MAP

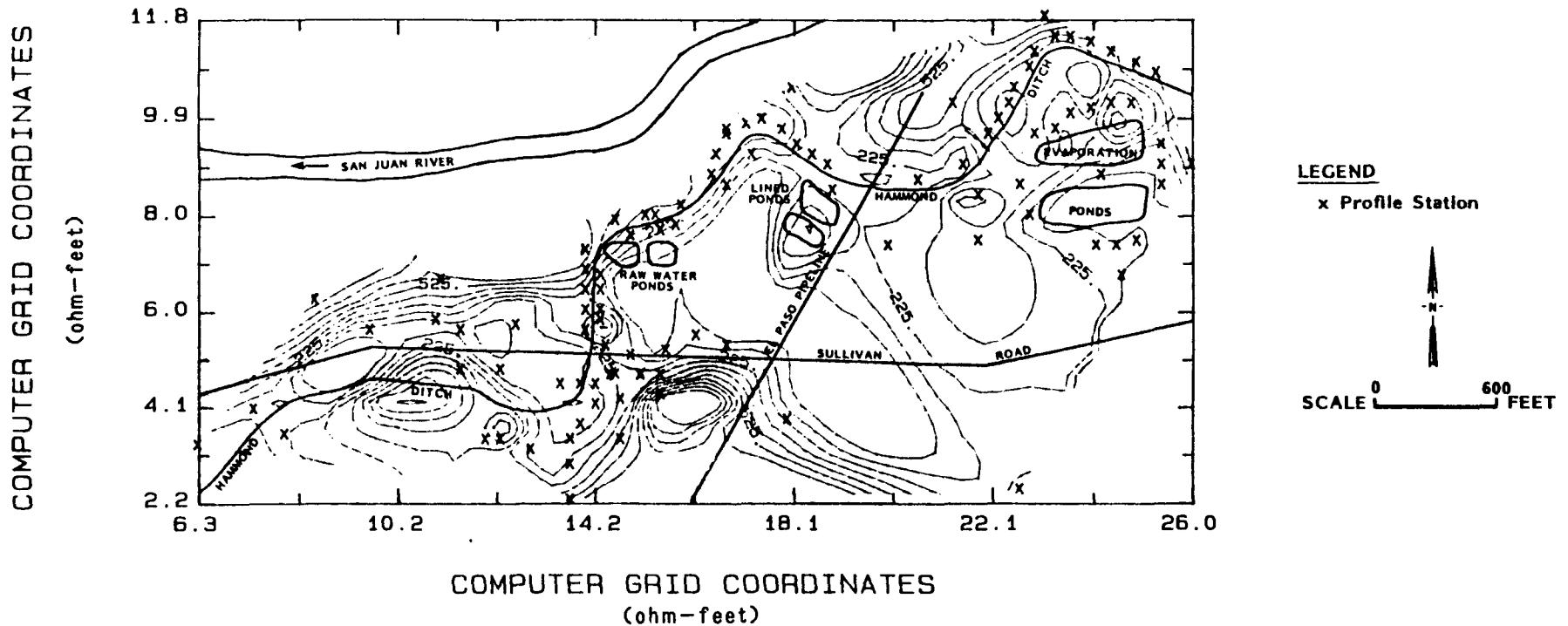


FIGURE 2.5

BLOOMFIELD REFINING CO.: 40-FT. ER PROFILE MAP

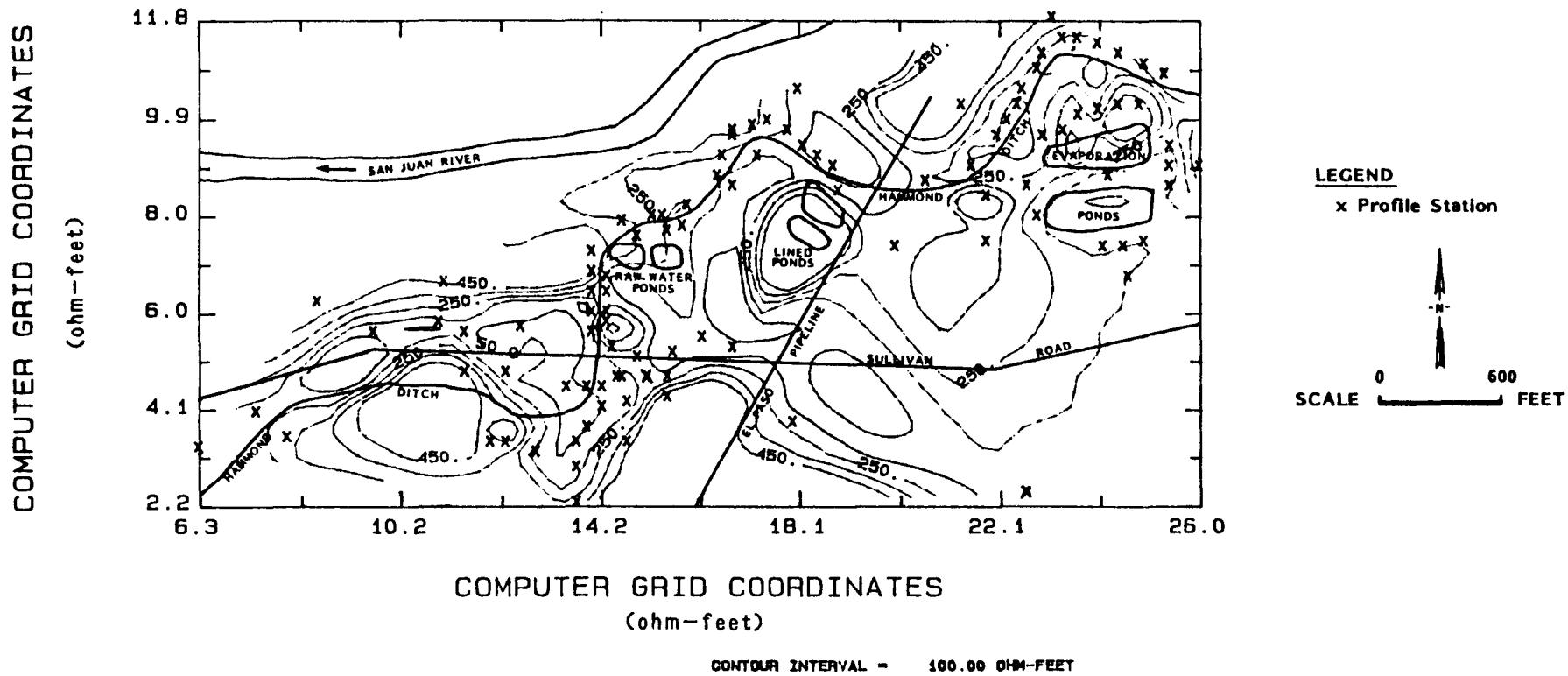


FIGURE 2.6

BLOOMFIELD REFINING CO.: 60-FT. ER PROFILE MAP

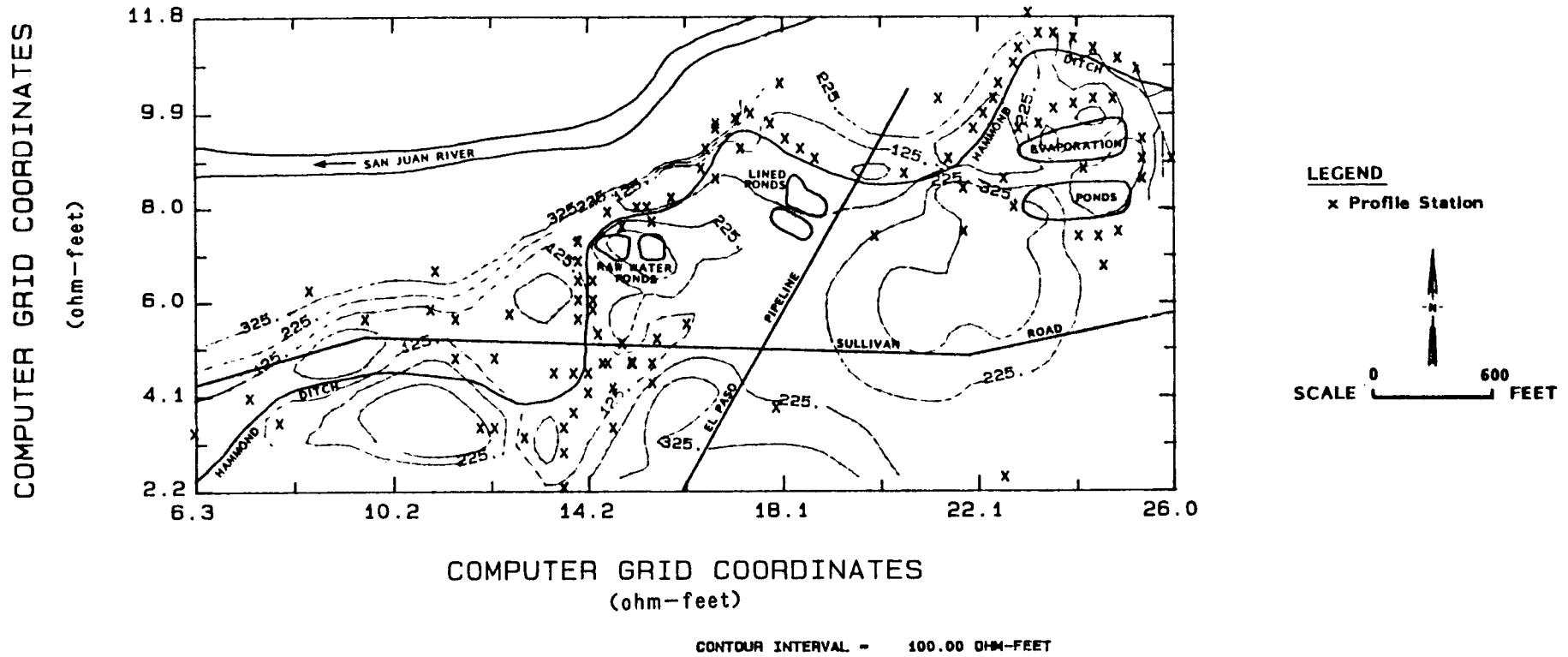


FIGURE 2.7

BLOOMFIELD REFINING CO.: 80-FT. ER PROFILE MAP

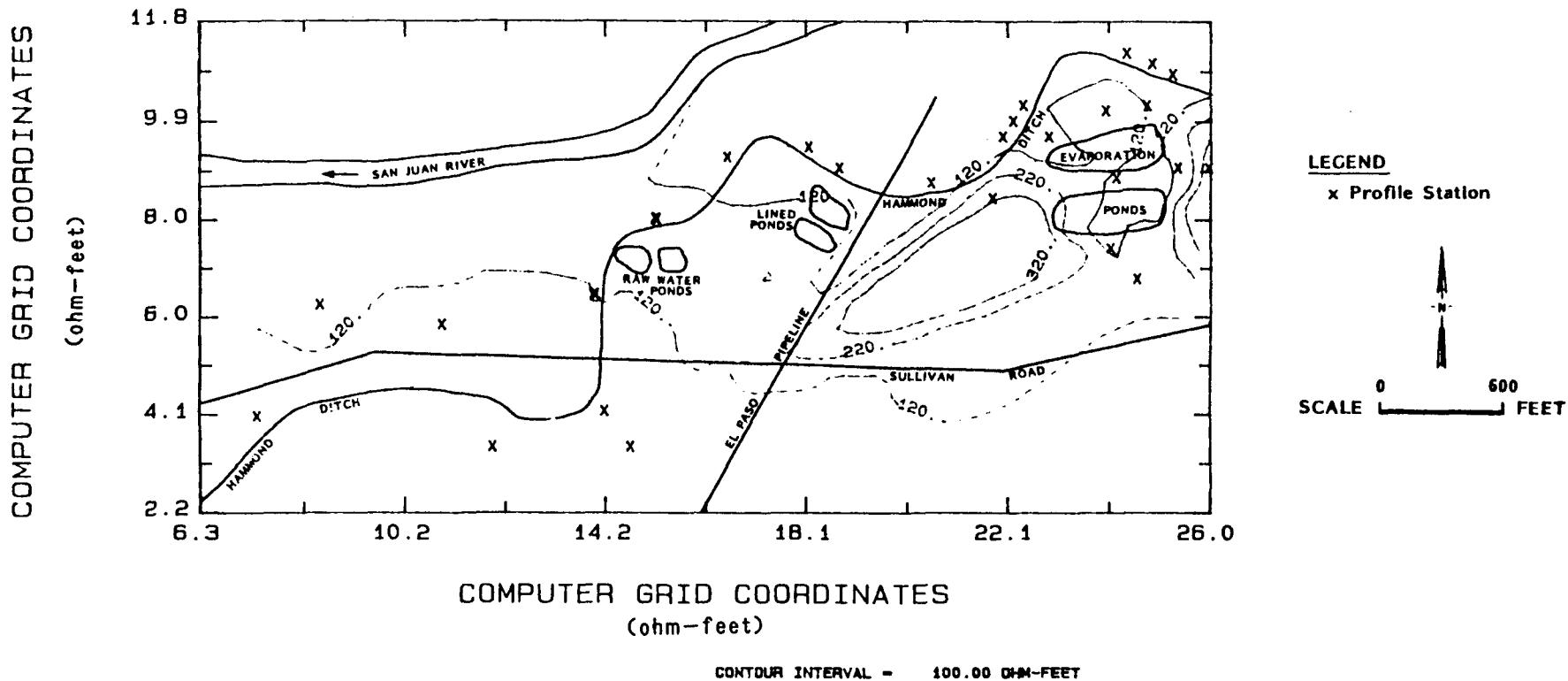


FIGURE 2.8

BLOOMFIELD REFINING CO.: 100-FT. ER PROFILE MAP

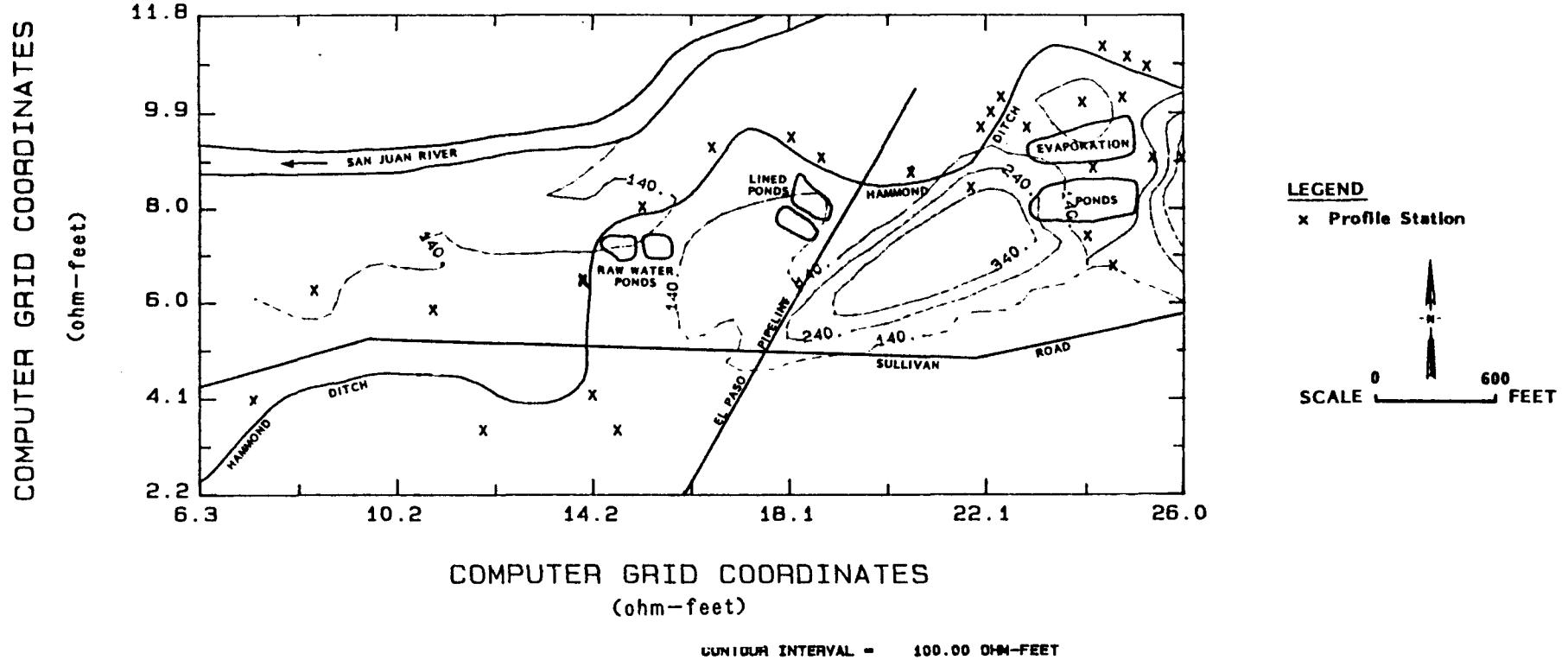


FIGURE 2.9
CROSS SECTION LOCATIONS

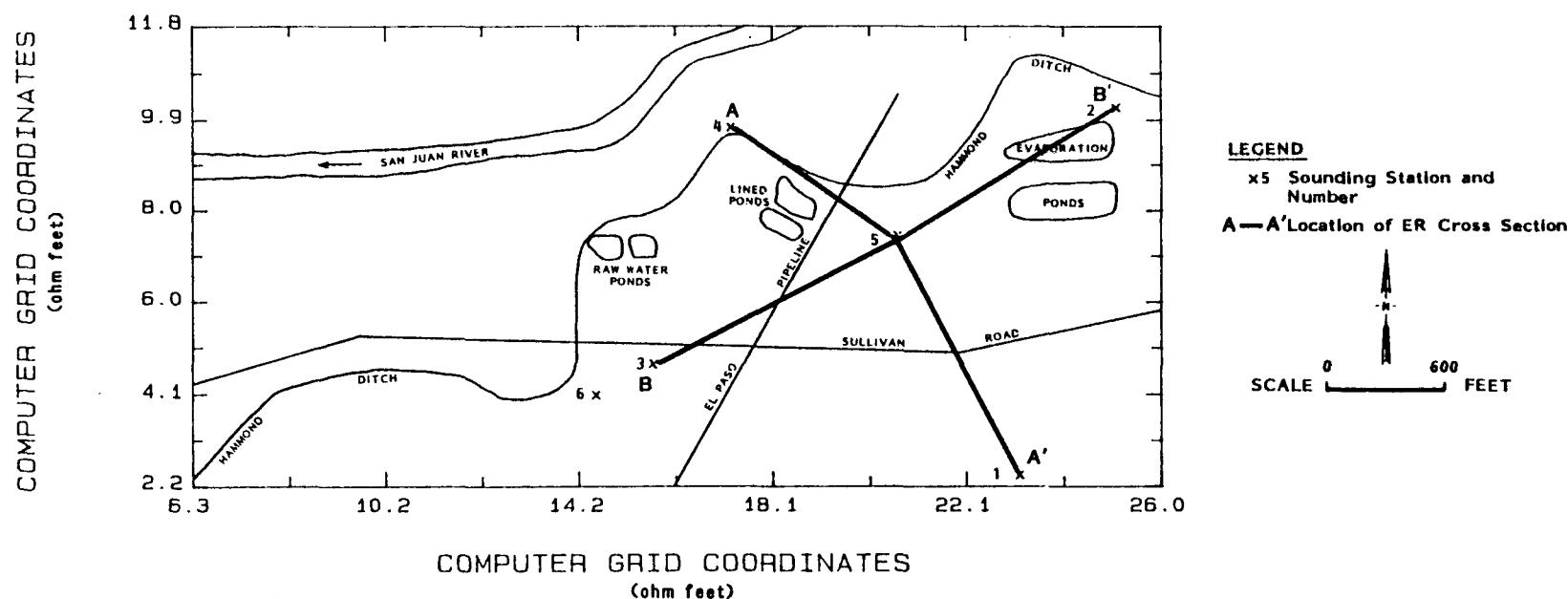


FIGURE 2.10

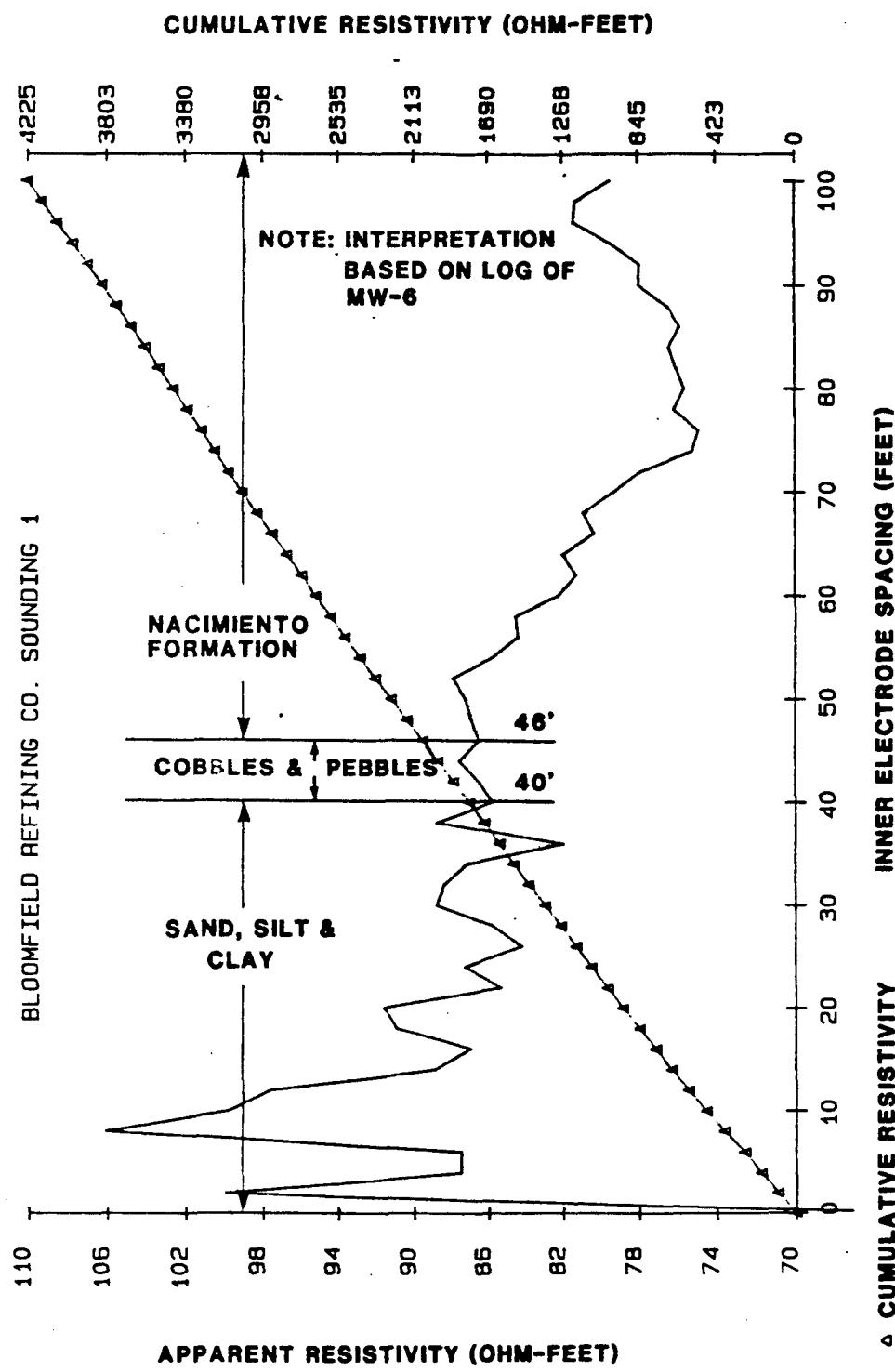


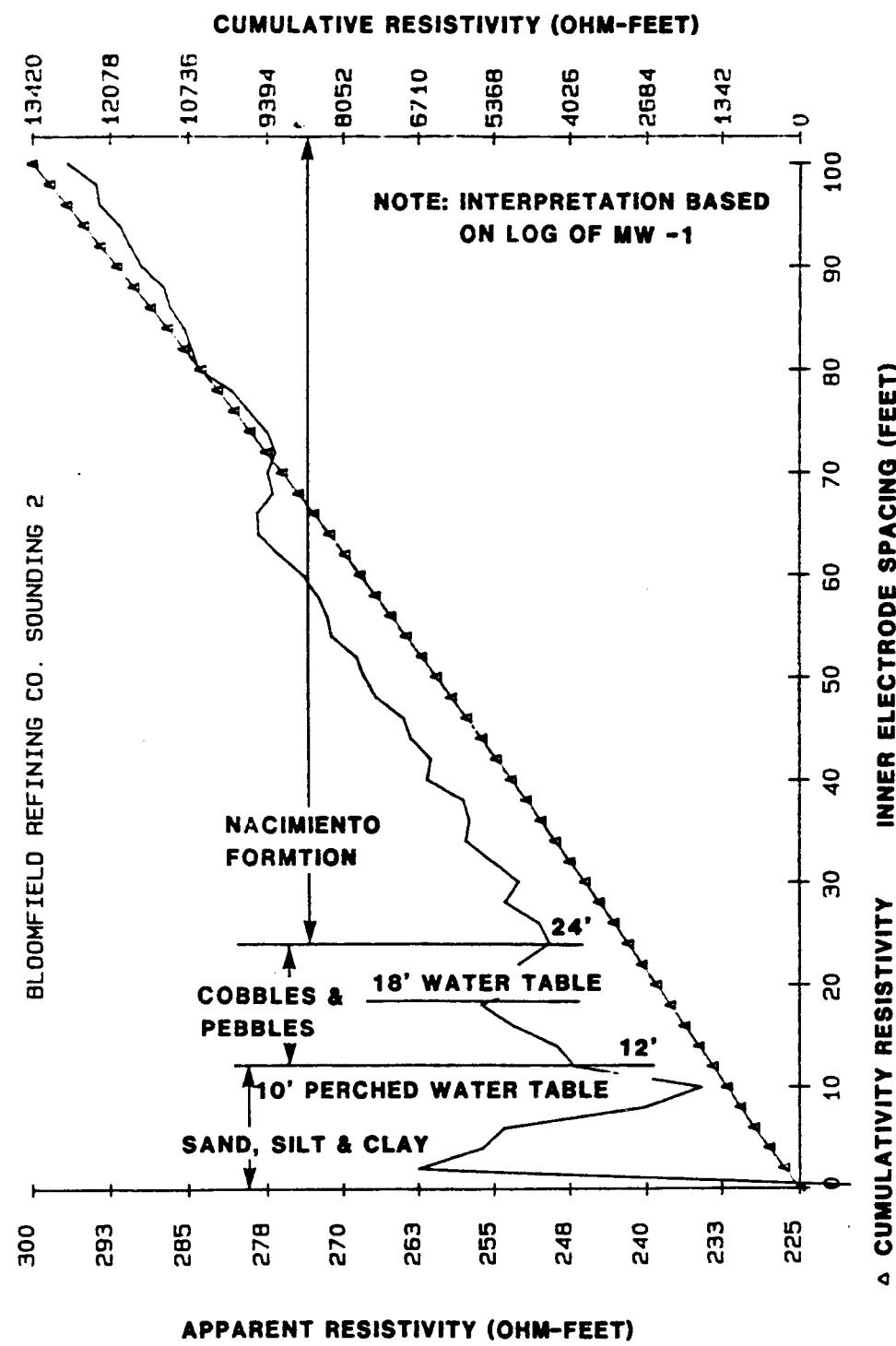
FIGURE 2.11

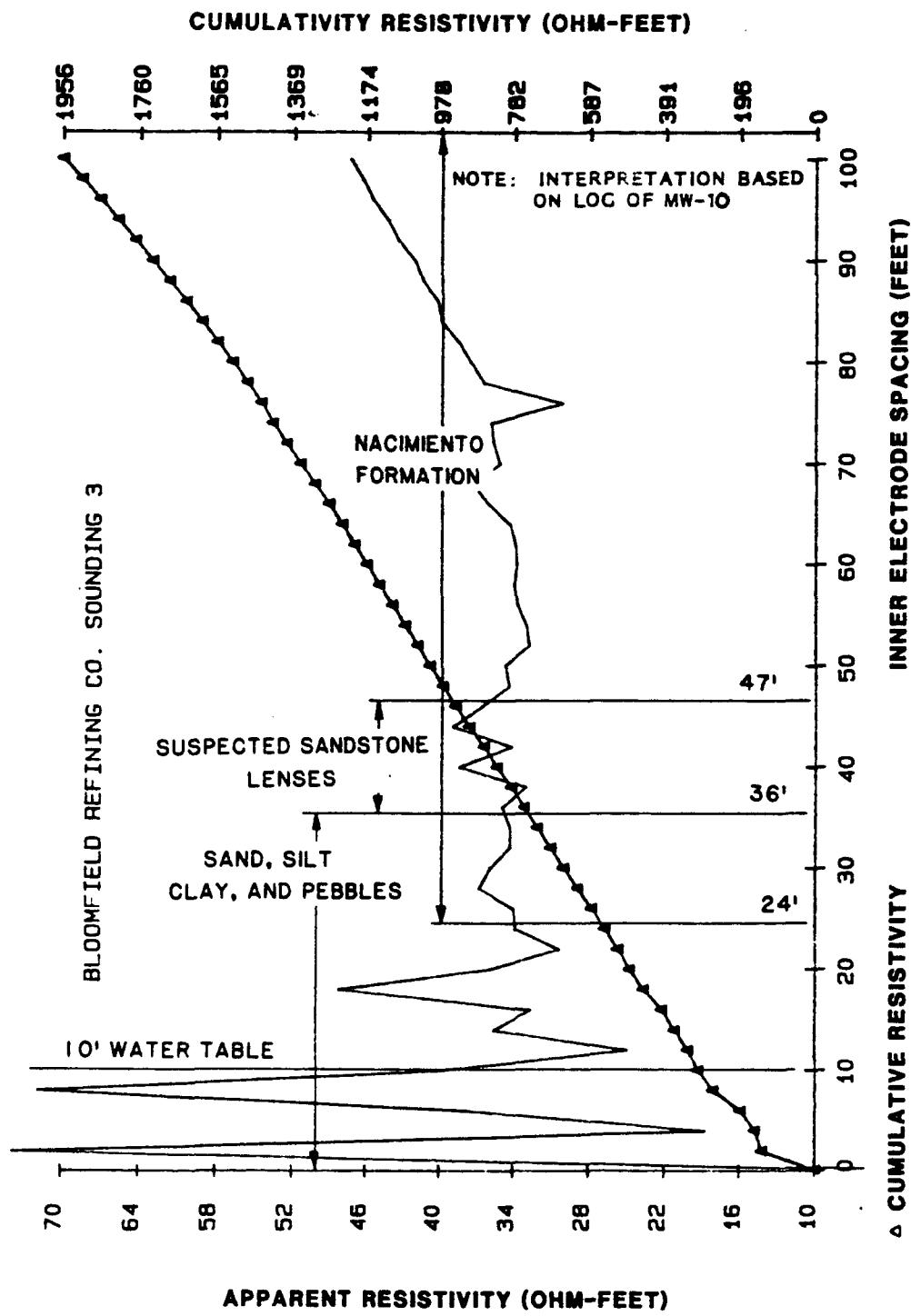
FIGURE 2.12

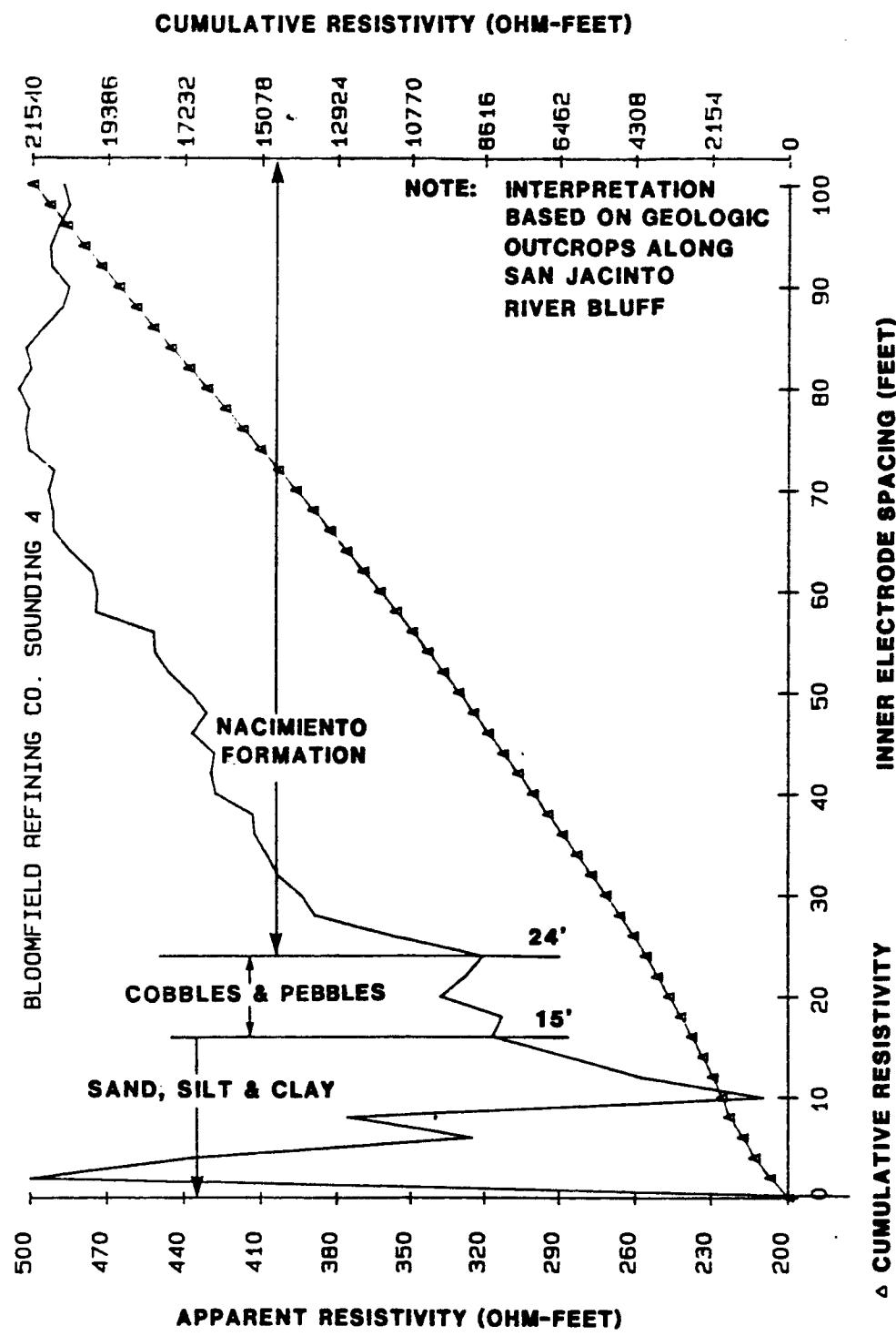
FIGURE 2.13

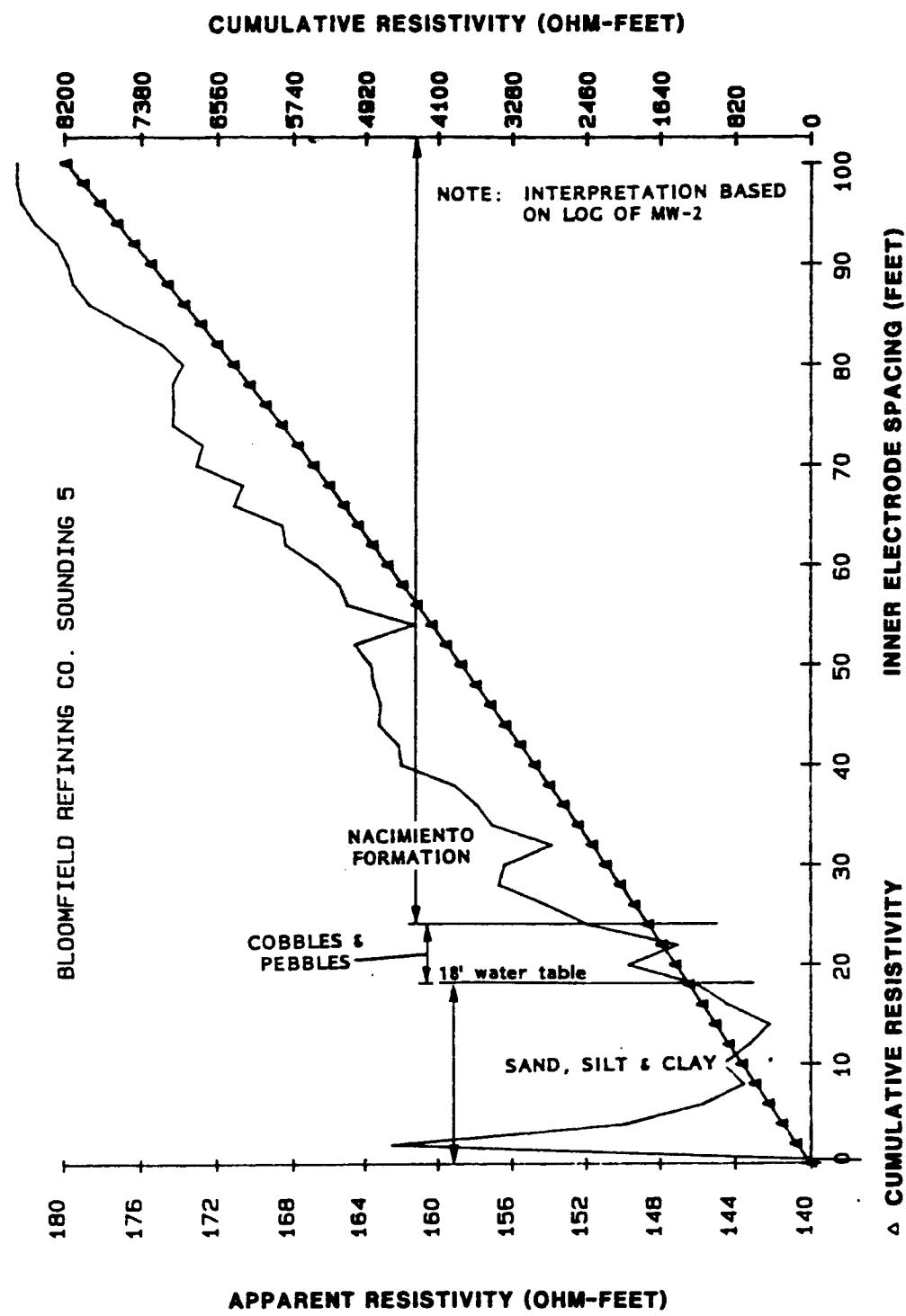
FIGURE 2.14

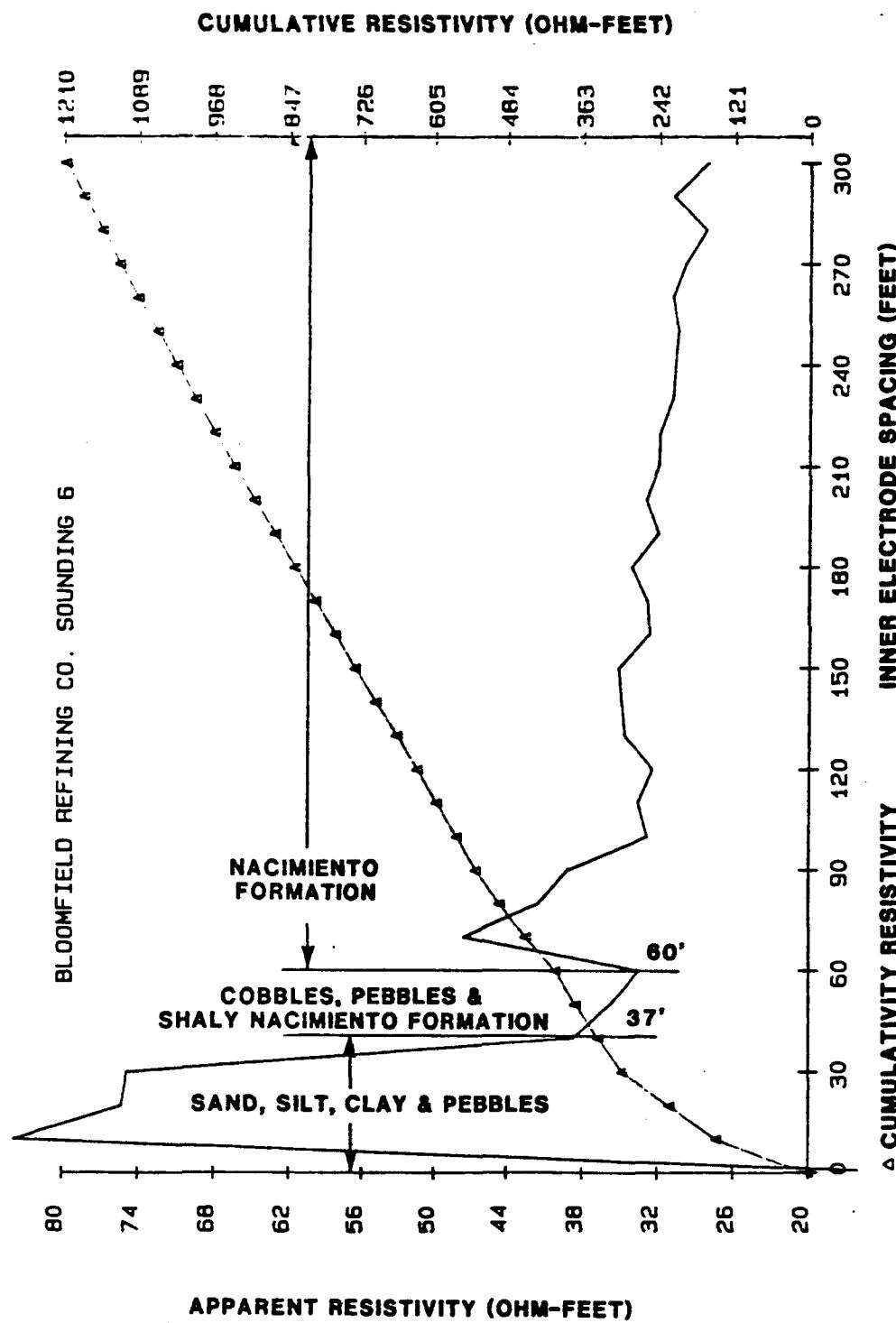
FIGURE 2.15

FIGURE 2.16

CROSS SECTION A-A'

A'

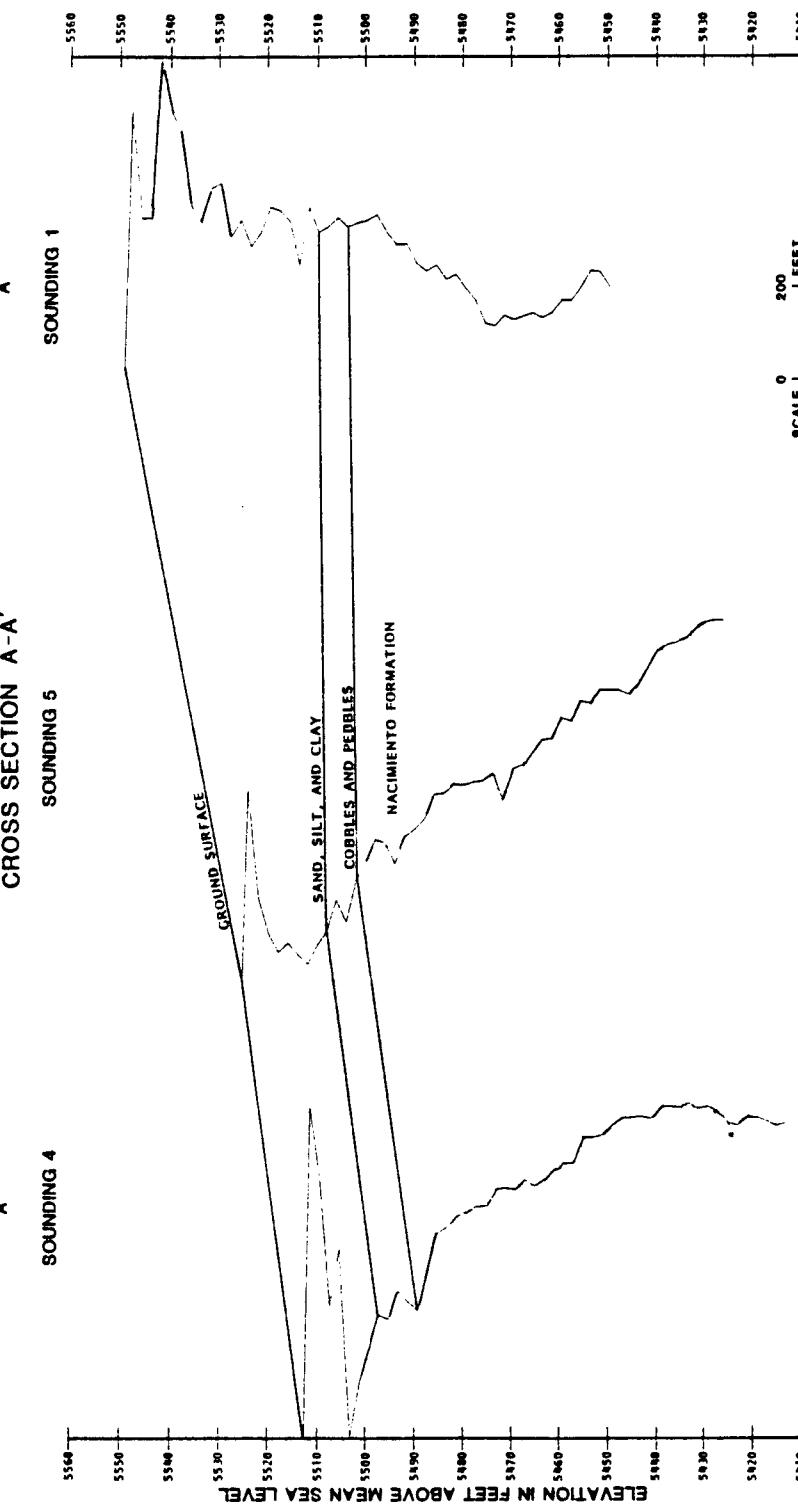


FIGURE 2.17
CROSS SECTION B-B'

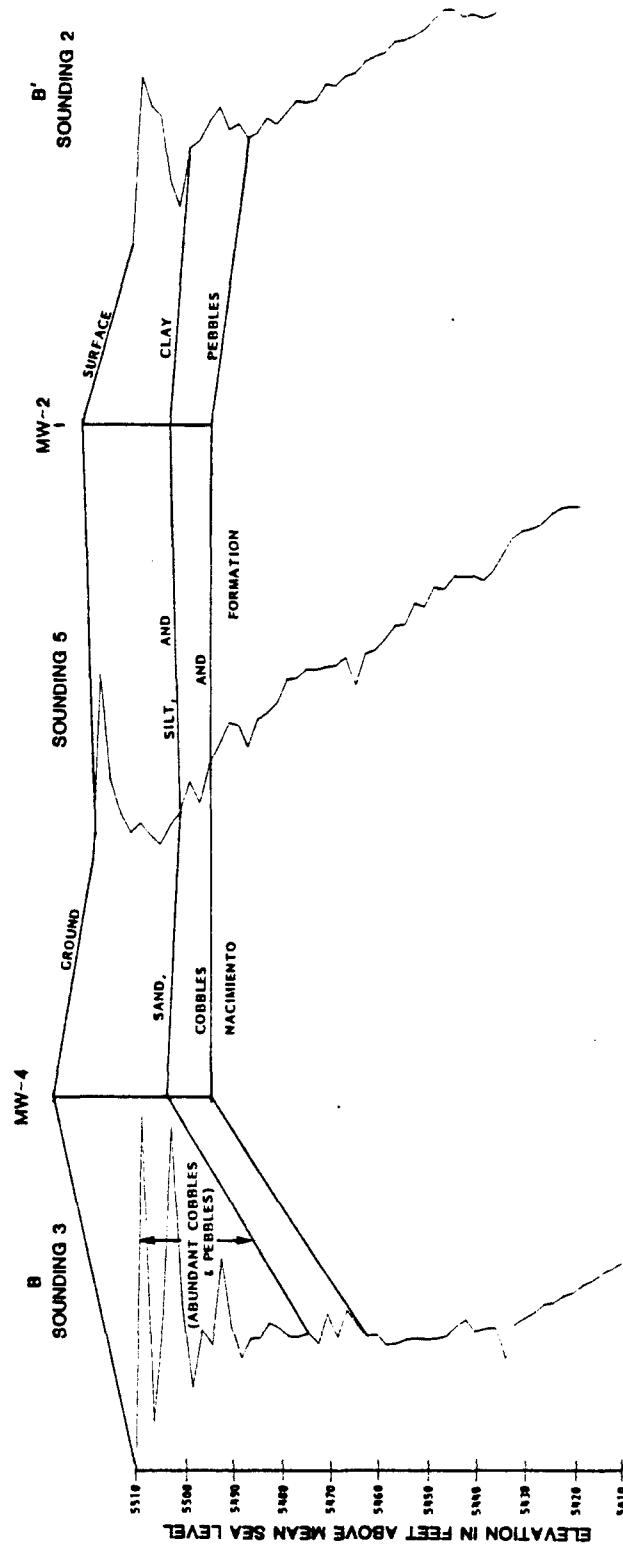
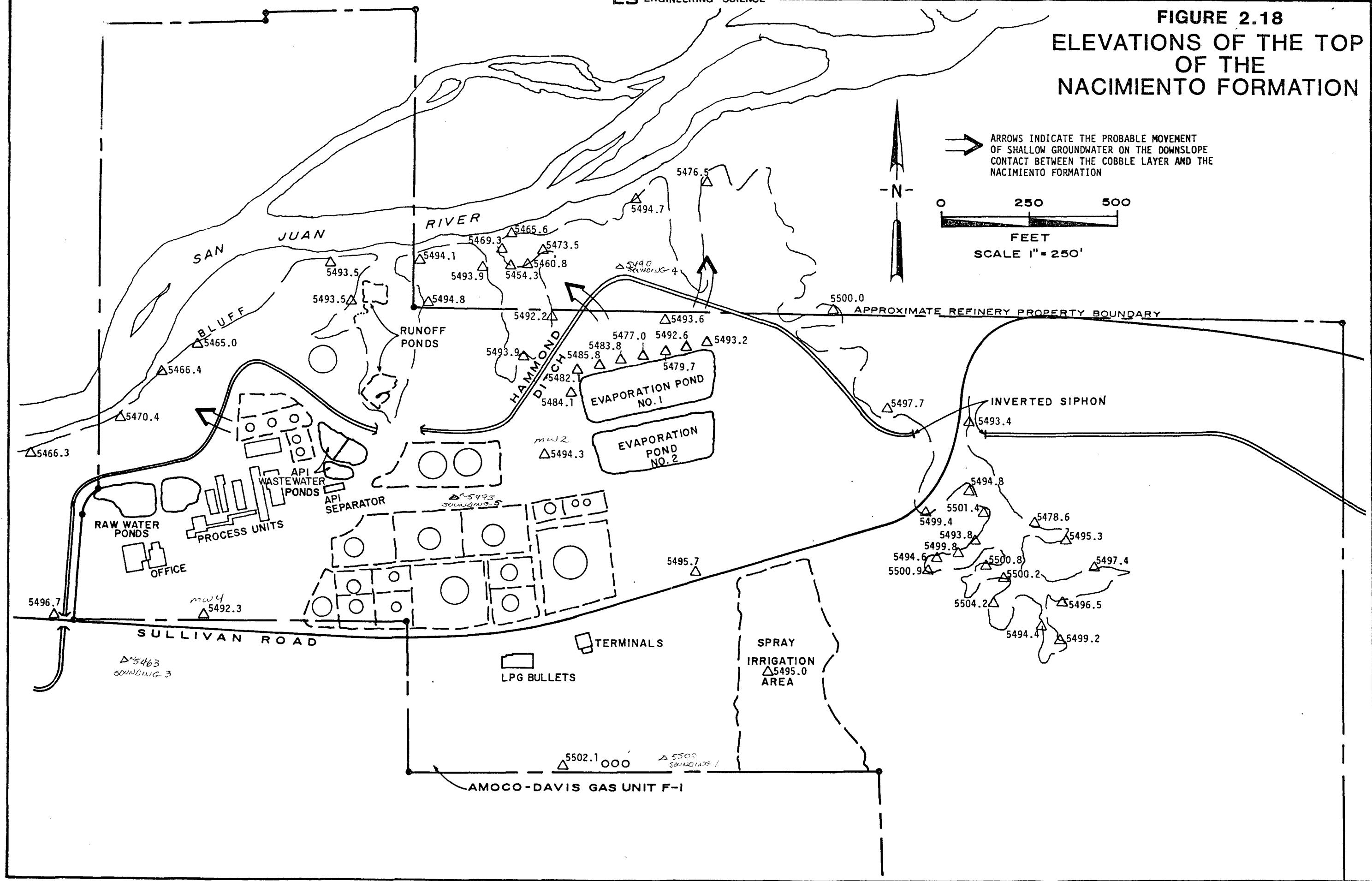


FIGURE 2.18
ELEVATIONS OF THE TOP
OF THE
NACIMIENTO FORMATION



New Mexico Oil Conservation District (NMOCD) for Plateau Refining by Dr. William Turner of American Groundwater Consultants, provides the elevations of the top of the Nacimiento. These data show the formation was subject to prehistoric erosion. This erosion created a series of subsurface channels which were filled with permeable materials by subsequent sedimentation processes. The net result of these geologic processes is the existence of a variable saturated zone thickness over the site which affects the route of groundwater movement.

DISCUSSION

The presence, areal extent, and route of migration of subsurface contamination can be interpreted from the ER data only in the context of how the ER data are either corroborated or refuted by other evidence. Since the data provide only a relative measure of resistance, the data by themselves can only suggest possible locations of subsurface hydrocarbons. Other ER signatures which can appear the same as hydrocarbon materials are changes in soil moisture, changes in soil or rock composition, or structural features. The ER data must therefore be "ground truthed" via some additional positive evidence.

The computer-generated ER profile maps should be evaluated with the understanding that the ohm-foot contours delineated on the maps constitute only an interpolation and extrapolation of the discrete ER measurements presented in Figure 2.1. The weight given to the significance of the contours should be tempered with an examination of Figure 2.1 to determine the density of ER data points in a particular area of the site.

Both the process area and refined product storage area provided little accessibility during the ER survey due to numerous pipeline interferences and structures. Therefore, the 25- to 50-ohm-foot contours representing this area are based on only a few data points.

Since relatively low ER values can suggest the presence of subsurface hydrocarbons, the areas of the BRC site showing relatively low ER

values are of interest. Based on the 10- to 40-foot depth profile maps, there are three areas showing relatively low ER values.

One area of relatively low ER contours (approximately 200 ohm-feet) exists to the west of the solar evaporation ponds and is generally oriented in a southeast to northwesterly direction and transects Hammond Ditch. In a discharge plan approved by NMOCD on June 7, 1984, the solar evaporation ponds were reported to be subject to some leakage and percolation (see references section). In addition, the area mentioned corresponds to an erosion channel in the Nacimiento (see Figure 2.18) which provides an avenue for flow from the ponds toward the San Juan River during nonirrigation periods in Hammond Ditch (see Chapter 5). The ER survey was conducted during the nonirrigation season. Thus, the low ER values may be the result of leakage of water from the ponds and not subsurface hydrocarbons.

Compare
with M.W.
results.

Another area depicted to show relatively low ER values is defined by the 25- to 50-ohm-foot contours on the 10- to 40-foot contour maps. This area includes both the process and refined product storage areas. As mentioned earlier, the exact extent of these low-valued contours is suspect because of the limited number of ER measurements in this area. Some indication of free product existed in monitoring well 4 during the ER survey, and this well was used as a frame of reference for other ER measurements in the area. However, monitoring wells 2, 3, and 5 through 10 are also screened within the same zones of low resistance included in the projected 25- to 50-foot contour areas and have not produced any indication of free product. Moreover, laboratory summaries of these wells' quarterly sampling results, shown in chapter 4, show low levels of hydrocarbon compounds.

The last location exhibiting low ER contours is off site in an area south and west of the intersection of Hammond Ditch and Sullivan Road. Based on the ER contours, these sites could represent the route of migration of groundwater measured in monitoring wells 4, 7, and 10. No significant investigation of this area is available at this time since the focus of the 3013 administrative order directives has been to

concentrate investigative efforts within the property boundaries of the refinery.

Recently, an independent assessment of offsite tracts has been initiated for the NMOCD. This study involves installation of a test recovery well in an area of low resistivity located on site near monitoring well 4. The study also involves installation of 2 monitoring wells in areas of similarly low resistivity located south and west of BRC's property. This study will examine recovery characteristics of the saturated zone and will monitor the saturated zone west of the property for subsurface hydrocarbons.

In summary, the ability to explain the existence of low ER contours is subject to procurement of additional information which is not available at this time. Monitoring data verify the low ER contours only in a localized area monitored by wells 4, 9, and 10. Available information developed from the administrative order work elements which establishes the basis for explanation of the low ER contours is provided in the following chapters.

Subsurface conditions in offsite areas have not been evaluated since investigation of these areas has been beyond the scope of work. It is recommended that investigation of the offsite areas be pursued under auspices of the NMOCD program.

CHAPTER 3

GROUNDWATER MONITORING WELLS

Four groundwater monitoring wells, designated 7 through 10, were installed consistent with the guidelines specified in the EPA-approved work plan. Figure 3.1 depicts the locations of these wells with respect to the pre-existing wells (1 through 6) and site features. The locations for these wells were selected in consultation with USEPA and state officials based on site information and geophysical information submitted previously. Installation of these wells took place during the period from February 26 to March 4, 1986. Boring logs prepared for each of these wells are presented in Tables 3.1 through 3.4. These logs provide both a geologic description of the strata penetrated at each well location and a description of the type of well hardware, packing, and seals installed.

Well 7 was installed to investigate possible contamination of the Nacimiento Formation by hydrocarbons. It was installed with a 10-foot stainless steel screen located in a gray to green clay at a depth range of approximately 50 to 60 feet. A sand filter pack was installed from the bottom of the boring to a depth of 45 feet. A 4-foot bentonite plug was installed on top of the sand pack, followed by a grout seal installed to the surface. The well is screened only within the Nacimiento Formation and is plugged to prevent vertical migration of hydrocarbons from above.

For completeness, lithologic logs for monitoring wells 1 through 6 have been added. These logs are as prepared by Earl & Sons, Inc. (the drillers) and are presented as Tables 3.5 to 3.10.

FIGURE 3.1

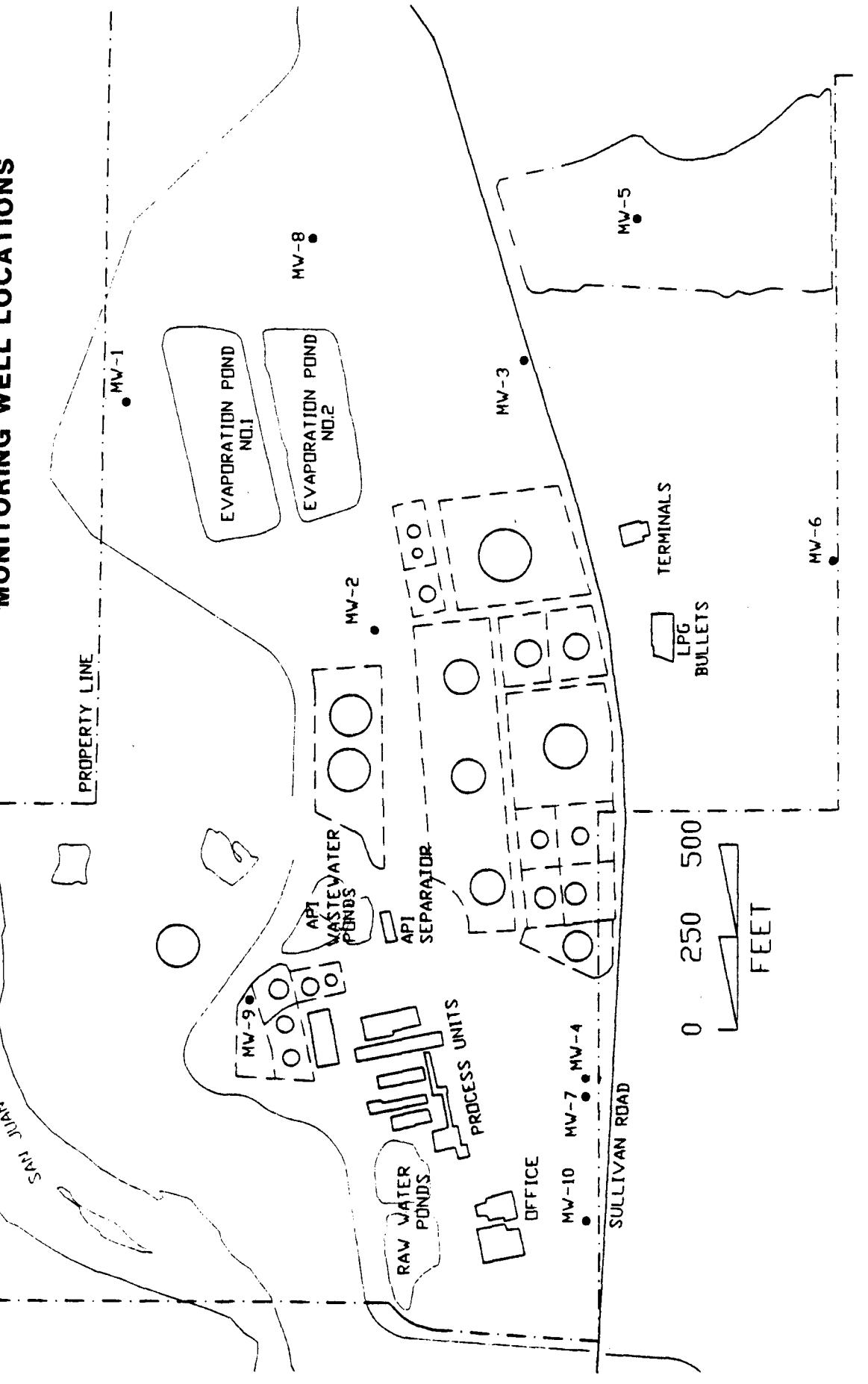
**BLOOMFIELD REFINERY
MONITORING WELL LOCATIONS**

TABLE 3.1
WELL LOG FOR MONITORING WELL NUMBER 7

Drilling Date: February 25

| <u>Depth in Feet</u> | <u>Description</u> |
|-----------------------------------|--|
| 0-1 | Gravel fill |
| 1-5 | Brown sandy silt and clay with small gravels |
| 5-10 | Brown sandy silt and clay, more firm and sticky |
| 10-15 | Lighter brown sandy silt and sticky clay |
| 15-20 | Lighter brown sandy silt and clay, larger cobbles and pebbles |
| 20-25 | Sand with cobbles and pebbles |
| 25-30 ^{26.5 ft. L.D.} | Sand |
| 30-35 | Greenish clay with pebbles, top of Nacimiento estimated at 32 feet |
| 35-40 | Greenish clay, few pebbles |
| 40-45 | Green to gray clay, smooth drilling |
| 45-50 | Green to gray clay, smooth drilling |
| 50-65 | Sticky gray to green clay |

Elevation of Top of Pipe: 5524.09 feet

Total Depth of Casing: 62.11 feet

Description of Casing: Bottom of casing has a 2 foot stainless steel blank section for a silt trap, followed by a 10 foot section of 6" I.D. stainless steel screen, in turn followed by 6" I.D. schedule 40 PVC casing to the top of pipe. Sand was added to 45 feet below grade, bentonite to 41 feet below grade, and grout to the surface.

Screen 50-60 ft
Sand pack 45-60 ft

Type n (L)
(likely Air)

TABLE 3.2
WELL LOG FOR MONITORING WELL NUMBER 8

Drilling Date: February 28, 1986

| <u>Depth in Feet</u> | <u>Description</u> |
|-----------------------------|---|
| 0-20 | Light brown sandy clay, similar to that found on the ground surface |
| 20-34 29 STRW | Cobbles and pebbles |
| 34 | Green-gray clay and sandstone, intermixed with small pebbles and sand. Top of Nacimiento. |

Elevation of Top of Casing: 5531.12 feet

Total Depth of Casing: 34.94 feet

Description of Casing: Bottom of casing has a 2 foot stainless steel blank section for a silt trap, followed by 20 feet of 6" I.D. stainless steel screen, followed by 6" I.D. schedule 40 PVC to the surface. The screened section of the hole was sanded to within 7 feet of the surface, a bentonite seal (1/2 bucket) was added and concrete was used for a surface seal.

*Screened 13 - 33 feet
Sand pack 7 - 33 feet*

TABLE 3.3
WELL LOG FOR MONITORING WELL NUMBER 9

Drilling Date: March 3, 1986

| <u>Depth in Feet</u> | <u>Description</u> |
|-----------------------------|---|
| 0-5 | Fill material, some rock |
| 5-10 | Sticky reddish brown silty clay |
| 10-15 | Lighter color silty clay, some pebbles |
| 15-20 | Lighter color silty clay, some pebbles |
| ^{20 feet} 20-25 | Cobbles, pebbles, sand |
| 25-30 | Cobbles, greenish clay, top of Nacimiento |

Elevation of Top of Casing: 5519.70 feet

Total Depth of Casing: 33.99 feet

Description of Casing: Bottom of casing has a 2 foot stainless steel blank section for a silt trap followed by 20 feet of 6" I.D. stainless steel screen, followed by 6" I.D. schedule 40 PVC to the surface. The screened section of the hole was sanded to within 7 feet of the surface, a bentonite seal (1/2 bucket) was added and concrete was used for a surface seal.

Screen 12-32
Sand 7-32

TABLE 3.4
WELL LOG FOR MONITORING WELL NUMBER 10

Drilling Date: March 4, 1986

| <u>Depth in Feet</u> | <u>Description</u> |
|--------------------------|--|
| 0-5 | Topsoil, roadbase, reddish brown sandy clay |
| 5-10 | Reddish brown silty, sandy clay |
| 10-15 | Cobbles, pebbles |
| 15-20 | Gravel, cobbles, pebbles |
| 20-25 | Greenish clay at 23 feet, top of Nacimiento |
| 25-30 | Greenish clay, Nacimiento |
| 30-35 | Nacimiento, color changed from yellow-green to blue-gray |

Elevation of Top of Casing: 5516.86 feet

Total Depth of Casing: 33.93 feet

Description of Casing: Bottom of casing has a 2 foot stainless steel blank section for a silt trap, followed by 20 feet of 6" I.D. stainless steel screen, followed by 6" I.D. schedule 40 PVC to the surface. The screened section of the hole was sanded to within 7 feet of the surface, a bentonite seal (1/2 bucket) was added and concrete was used for a surface seal.

Screen 12-32
Sand 7-32

qc of #1D completed in
clay & Nacimiento.
11' in alluvium.
Sand above top of
screen.

TABLE 3.5
WELL LOG FOR MONITORING WELL NUMBER 1

Drilling Date: February 8, 1984
Location: 29.11.27.24221

| <u>Depth in Feet</u> | <u>Description</u> |
|---------------------------|--|
| 0-5 | Light brown clayey sand, coarse, poorly sorted, quartzose, and slightly calcareous |
| 5-10 | Yellowish gray sandy pebbles and cobbles, poorly sorted, rounded to subrounded |
| 10-12 | Yellowish gray pebbly sand, very coarse, poorly sorted, felospathic and noncalcareous |
| 12-22 <i>16.7 ft w</i> | Dark gray pebbly and sandy cobbles, some quartz pebbles, most are volcanic, subrounded cobbles and pebbles, some clay, a little water at about 16 feet |
| 22-25 | Gray-green clayey sand becoming light yellow clayey sandstone and sandy claystone |

TABLE 3.6
WELL LOG FOR MONITORING WELL NUMBER 2

Drilling Date: February 7, 1984
Location: 29.11.27.24321

| <u>Depth in Feet</u> | <u>Description</u> |
|------------------------------|--|
| 0-5 | Light yellow brown silty sandy clay, very calcareous |
| 5-10 | Light yellow brown clayey sand, subrounded to subangular, moderately to poorly sorted, very calcareous |
| 10-15 | Light brown pebbly sand, clayey, very calcareous, cobbles at 15 feet |
| 15-20 <i>18.5 L T. S.</i> | Gray sandy pebbles, poorly sorted coarse quartzose sand, pebbles are dark gray and volcanic |
| 20-25 | Dary gray cobbles, some quartz pebbles, mostly volcanic, some sand |
| 25-26 | Yellow gray clayey sandstone and sandy claystone |

TABLE 3.7
WELL LOG FOR MONITORING WELL NUMBER 3

Drilling Date: February 8, 1984
Location: 29.11.27.24443

| <u>Depth in Feet</u> | <u>Description</u> |
|----------------------------------|--|
| 0-5 | Yellow brown sandy silt and clay, very calcareous quartzose |
| 5-10 | Yellow brown sand, calcareous, silty and clayey, quartzose |
| 10-15 | Yellow brown sand, silty and clayey, fine-grained, very calcareous, quartzose |
| 15-27 | Light brown clay, sandy, very clacareous, becoming pebbly with depth |
| 27-35 <small>33.45 ft</small> | Gray yellow brown cobbly sand, coarse, poorly sorted, silty and clayey, volcanic pebbles, small amount of water at about 35 feet |
| 35-40 | Gray cobbles, pebbly and sandy, coarse sand, yellow gray clayey sandstone at about 40 feet |

TABLE 3.8
WELL LOG FOR MONITORING WELL NUMBER 4

Drilling Date: February 9, 1984
Location: 29.11.27.23344

| <u>Depth in Feet</u> | <u>Description</u> |
|--------------------------|--|
| 0-5 | Yellow gray-brown sandy silt and clay, calcareous |
| 5-10 | Yellow brown silty sandy clay and clayey silt, very slightly calcareous |
| 10-15 | Reddish yellow-brown clayey sandy silt, silty clay, fine-grained quartzose sand, noncalcareous |
| 15-19 | Light brown coarse sand with clay and pebbles, calcareous |
| 19-25 | Gray pebbly sand, very coarse, poorly sorted, some clay and silt, subrounded to subangular, quartzose, pebbles rounded, slightly calcareous |
| 25-30 | Gray cobbles and pebbles, subrounded to rounded, volcanic; at about 28 feet, <u>hydrocarbon smell and color</u> |
| 30-32 | Gray cobbly sand, with hydrocarbon smell and color, coarse-grained, sand is quartzose and feldspathic, subrounded and subangular quartz grains are clear |
| 32 | Yellow gray clayey sandstone |

TABLE 3.9
WELL LOG FOR MONITORING WELL NUMBER 5

Drilling Date: February 6, 1984
Location: 29.11.26.31112

| <u>Depth in Feet</u> | <u>Description</u> |
|--------------------------|--|
| 0-5 | Pale yellow brown clay, silty, some sand, calcareous |
| 5-10 | Pale yellow brown clayey sand and quartzose silt, poorly sorted, calcareous |
| 10-15 | Yellow brown sand, subrounded quartzose sand slightly calcareous |
| 15-20 | Yellow brown sand, clayey, moderately coarse-grained, very calcareous |
| 20-25 | Yellow brown sand, clayey, silty, fine- to medium-grained, moderately sorted, noncalcareous |
| 25-35 | Yellow brown sand, silty and slightly clayey, fine- to medium-grained, well sorted, subangular, noncalcareous, becoming more clayey with depth |
| 35-37 | Yellow brown pebbly and cobbly sand, clayey, calcareous |
| 37-47 | Dark gray sandy and clayey cobbles and pebbles, water at 42 feet |
| 47-50 | Dark gray cobbles with greenish clay |
| 50-54 | Green-gray pebbly clay |

TABLE 3.10
WELL LOG FOR MONITORING WELL NUMBER 6 *DRY*

Drilling Date: February 7, 1984
Location: 29.11.27.42144 or 42233

| <u>Depth in Feet</u> | <u>Description</u> |
|--------------------------|--|
| 0-15 | Pale yellow brown sand, clayey and silty, subangular, poorly sorted, quartzose, very calcareous, becoming more clayey with depth |
| 15-20 | Pale yellow brown silt, sandy and clayey, silt is coarse, sand is very fine, moderate sorting, quartzose and calcareous |
| 20-25 | Pale yellow sand, slightly clayey, subrounded, well sorted, quartzose, noncalcareous |
| 25-35 | Pale yellow sand, coarse- to medium-grained, quartzose, noncalcareous |
| 35-41 | Pale yellow sand, clayey, fine-grained, silty, quartzose, slightly calcareous |
| 41-49 | Gray-black cobbles and pebbles, volcanic |
| 49-52 | Gray-green clayey sandstone and sandy claystone |

CHAPTER 4

GROUNDWATER MONITORING WELL SAMPLING

BACKGROUND

Based on the requirements of the approved work plan, each of the groundwater monitoring wells at the Bloomfield refinery has been sampled by Bloomfield staff quarterly over a period of 12 months. Water samples from wells 1 through 3 and 5 were analyzed for priority pollutant heavy metals, cyanide, phenols, TOC, TDS, chloride, sulfate, and volatile organics including benzene, toluene, xylene, and ethylbenzene. In addition, at monitoring well 1, selected organic priority pollutants (see Appendix A) were analyzed during the second and fourth quarters. Water samples from monitoring wells 4 and 7 through 10 were analyzed for base/neutral priority pollutants in addition to the constituents already listed. Monitoring well 6 has been dry and therefore has never yielded any water for analyses.

FINDINGS

The laboratory results corresponding to each of the quarterly sampling efforts are presented in Appendix A to this report. In addition, sampling results for Hammond Ditch (covered in chapter 7) are presented in Appendix B to this report. Summaries of these data provide concentrations of detectable substances and laboratory detection limits. Field measurements of pH, conductivity, and water levels are also included. These summaries give a concise picture of each well's performance through the sampling period in Tables 4.1 through 4.9. As a reference, Appendix C presents EPA ambient standards and criteria for superfund remedial sites.

DISCUSSION

nd A review of the summaries shows that monitoring wells 1, 4, 7, 8, and 10 all show appearance of low levels of priority pollutants in only a single quarter's sampling effort. The lack of consistency in appearance

TABLE 4.1
RCRA 3013
GROUNDWATER RESULTS SUMMARY*
MW-1

| Parameter | 3/26/86 | 6/23/86 | 9/18/86 | 12/16/86 | Nominal Detection Limits |
|----------------------------|---------|---------|---------|----------|--------------------------|
| Laboratory Analysis (mg/l) | | | | | |
| Cyanide | | 0.1 | 0.07 | | 0.01 |
| Total phenols | 0.009 | 0.017 | 0.19 | 0.012 | 0.001 |
| TOC | 18 | 24 | 24 | 18 | 0.1 |
| TDS | 2,936 | 2,960 | 2,866 | 2,498 | 1 |
| Chloride | 750** | 994.7 | 814 | 774 | 1 |
| Sulfate | 7.5** | 630 | 673 | 579 | 1 |
| Antimony | | | | 0.25 | 0.01 |
| Arsenic | | 0.077 | 0.05 | | 0.05 |
| Beryllium | | | | 0.02 | 0.01 |
| Cadmium | 0.050 | | | | 0.01 |
| Lead | 0.085 | 0.065 | 0.15 | | 0.05 |
| Nickel | 0.08 | | 0.07 | 0.06 | 0.06 |
| Selenium | | 0.035 | 0.033 | 0.03 | 0.01 |
| Zinc | | 0.20 | 0.04 | 0.012 | 0.01 |
| Barium | NA | | | 0.055 | 0.005 |
| Manganese | NA | 0.25 | | 1.11 | 0.005 |
| Aluminum | NA | 2.07 | | 4.54 | 0.05 |
| Boron | NA | | | 0.27 | 0.01 |
| Molybdenum | NA | | | 0.17 | 0.01 |
| Nitrate-nitrogen | NA | 0.54 | | 2.9 | 0.1 |
| 1,2-dichloroethane | | | | 0.002 | 0.001 |
| Physical Measurements | | | | | |
| pH, field, s.u. | 7.30 | 7.25 | 7.27 | 7.19 | |
| Conductivity | | 4600 | 4600 | 4400 | |
| T.O.C. elevation, ft | 5515.77 | 5515.77 | 5515.77 | 5515.77 | |

NA - Not analyzed for that date.

Since heavy metals were not filtered, could have particulates dissolved in acid.
Lower SO₄, High TOC

TABLE 4.1 (Continued)

RCRA 3013
GROUNDWATER RESULTS SUMMARY¹
MW-1

| Parameter | 3/26/86 | 6/23/86 | 9/18/86 | 12/16/86 | Nominal Detection Limits |
|------------------------------------|---------|---------|---------|----------|--------------------------------|
| Depth to groundwater, ft | 16.70 | 14.56 | 15.74 | 16.32 | |
| Elevation of groundwater, ft | 5499.07 | 5501.21 | 5500.03 | 5499.45 | |

*Summary includes only pollutants determined to be present at concentrations greater than detection limits.

**Laboratory technique error detected.

TABLE 4.2
RCRA 3013
GROUNDWATER RESULTS SUMMARY*
MW-2

| Parameter | 3/26/86 | 6/23/86 | 9/18/86 | 12/16/86 | Nominal Detection Limits |
|-----------------------------------|---------|---------|---------|----------|--------------------------|
| Laboratory Analysis (mg/l) | | | | | |
| Cyanide | | 0.1 | 0.18 | | 0.01 |
| - Total phenols | 0.063 | 0.023 | 0.17 | 0.110 | 0.001 |
| TOC | 18 | 27 | 23 | 15 | 0.1 |
| TDS | 2,796 | 3,650 | 3,598 | 3,664 | 1 |
| - Chloride | 200** | 1,204.6 | 993 | 1,012 | 1 |
| - Sulfate | 11.0** | 1,750 | 1,104 | 1,372 | 1 |
| Antimony | | | | 0.48 | 0.01 |
| Arsenic | | 0.094 | 0.08 | | 0.05 |
| Lead | 0.12 | | 0.08 | | 0.05 |
| Nickel | 0.07 | | 0.12 | 0.08 | 0.06 |
| - Selenium | | 0.070 | 0.104 | 0.04 | 0.01 |
| Silver | 0.003 | | | | 0.002 |
| Zinc | | 0.020 | 0.02 | 0.009 | 0.01 |
| Physical Measurements | | | | | |
| pH, field, s.u. | 7.23 | 7.17 | 6.78 | 7.22 | |
| Conductivity | | 5400 | 5500 | 5800 | |
| T.O.C. elevation, ft | 5519.45 | 5519.45 | 5519.45 | 5519.45 | |
| Depth to groundwater, ft | 18.80 | 18.27 | 18.23 | 18.4 | |
| Elevation of groundwater, ft | 5500.76 | 5501.18 | 5501.22 | 5501.05 | |

*Summary includes only pollutants determined to be present at concentrations greater than detection limits.

**Laboratory technique error detected.

TABLE 4.3
RCRA 3013
GROUNDWATER RESULTS SUMMARY*
MW-3

| Parameter | 3/26/86 | 6/23/86 | 9/18/86 | 12/16/86 | Nominal Detection Limits |
|-----------------------------------|---------|---------|---------|----------|--------------------------|
| Laboratory Analysis (mg/l) | | | | | |
| - Cyanide | | 0.25 | 0.17 | 0.07 | 0.1 |
| - Total phenols | 0.006 | 0.006 | 0.082 | 0.012 | 0.001 |
| TOC | 29 | 17 | 16 | 12 | 0.1 |
| - TDS | 4,836 | 5,362 | 5,514 | 4,860 | 1 |
| Chloride | 1,500** | 1,584 | 1,290 | 1,290 | 1 |
| - Sulfate | 29.5** | 1,950 | 2,056 | 2,204 | 1 |
| Toluene | | 0.003 | | | 0.001 |
| Xylene | | 0.030 | | | 0.001 |
| Antimony | | | | 0.67 | 0.01 |
| - Arsenic | | 0.15 | 0.21 | | 0.05 |
| - Cadmium | 0.12 | 0.015 | | 0.11 | 0.01 |
| Lead | 0.14 | 0.070 | 0.18 | | 0.05 |
| - Mercury | 0.004 | | | | 0.002 |
| Nickel | 0.08 | 0.08 | 0.14 | 0.10 | 0.06 |
| - Selenium | | 0.010 | 0.100 | 0.05 | 0.01 |
| Zinc | | 0.018 | 0.018 | 0.01 | 0.01 |
| Physical Measurements | | | | | |
| pH, field, s.u. | 7.08 | 7.10 | 7.06 | 7.12 | |
| Conductivity | | 6900 | 7200 | 6900 | |
| T.O.C. elevation, ft | 5535.85 | 5535.85 | 5535.85 | 5535.85 | |
| Depth to groundwater, ft | 32.94 | 32.80 | 33.08 | 33.05 | |
| Elevation of groundwater, ft | 5502.91 | 5503.05 | 5502.77 | 5502.8 | |

*Summary includes only pollutants determined to be present at concentrations greater than detection limits.

**Laboratory technique error detected.

TABLE 4.4
RCRA 3013
GROUNDWATER RESULTS SUMMARY*
MW-4

| Parameter | 3/26/86 | 6/23/86 | 9/18/86 | 12/16/86 | Nominal Detection Limits |
|-----------------------------------|---------|---------|---------|----------|--------------------------|
| Laboratory Analysis (mg/l) | | | | | |
| - Cyanide | | 0.5 | | | 0.01 |
| - Total phenols | 0.633 | 0.430 | 0.085 | 0.096 | 0.001 |
| TOC | 110 | 130 | 63 | 170 | 0.1 |
| - TDS | 1,868 | 2,266 | 2,398 | 2,128 | 1 |
| - Chloride | 500** | 989.7 | 754 | 675 | 1 |
| Sulfate | | 12.5 | | | 1 |
| - Benzene | 11.8 | 3.1 | 6.65 | 1.91 | 0.001 |
| - Toluene | 7.5 | 0.290 | 0.407 | 1.78 | 0.001 |
| - Ethylbenzene | 0.107 | 0.070 | 0.140 | 4.48 | 0.001 |
| Antimony | | | | 0.40 | 0.01 |
| Arsenic | | 0.070 | 0.08 | | 0.05 |
| - Cadmium | 0.060 | | | | 0.01 |
| - Lead | 0.074 | 0.066 | | | 0.05 |
| Nickel | 0.08 | | 0.12 | | 0.06 |
| - Selenium | | 0.080 | 0.063 | 0.03 | 0.01 |
| Zinc | | 0.019 | 0.008 | 0.04 | 0.01 |
| - Barium | NA | 3.54 | | 2.3 | 0.005 |
| - Iron | | 12.0 | | 18.6 | 0.3 |
| - Manganese | | 3.5 | | 5.7 | 0.005 |
| Aluminum | | 1.93 | | 3.8 | 0.05 |
| Boron | | | | 0.7 | 0.01 |
| Fluoride | | 0.21 | | | 0.01 |
| Nitrate-nitrogen | | | | 0.41 | 0.1 |
| - 2,4-Dichloro-phenol | 0.200 | | | | 0.001 |
| - 2,4-Dimethyl-phenol | | 0.058 | | | 0.001 |
| - 4,6-Dinitro-o-cresol | 0.100 | | | | 0.001 |
| - 2,4-Dinitro-phenol | 0.050 | | | | 0.001 |
| - 2-Nitrophenol | | 0.108 | 0.026 | | 0.001 |
| - 4-Nitrophenol | 0.090 | 0.302 | 0.331 | | 0.001 |

Low SO₄, N-High TOC
possible reduction by
bacteria (See H.m)
P.116.

TABLE 4.4 (Continued)

RCRA 3013
GROUNDWATER RESULTS SUMMARY*
MW-4

| Parameter | 3/26/86 | 6/23/86 | 9/18/86 | 12/16/86 | Nominal Detection Limits |
|------------------------------|---------|---------|---------|----------|--------------------------|
| Phenol | 0.202 | | | | 0.001 |
| Benzo(a)anthra- | | | | | |
| cene | | 0.016 | 0.010 | | 0.001 |
| Chrysene | | 0.012 | | | 0.001 |
| Fluorene | 0.150 | | | 0.023 | 0.001 |
| Naphthalene | 0.036 | 0.019 | 0.015 | 0.036 | 0.001 |
| Pyrene | 0.166 | | 0.005 | | 0.001 |
| 2-Chloro-phenol | | | 0.001 | | 0.001 |
| P-chloro-m-cresol | | | 0.045 | | 0.001 |
| Acenaphthene | 0.044 | | | 0.049 | 0.001 |
| Physical Measurements | | | | | |
| pH, field, s.u. | 6.84 | 6.85 | 6.70 | 6.73 | |
| Conductivity | | 3800 | 3900 | 3800 | |
| T.O.C. elevation, ft | 5524.30 | 5524.30 | 5524.30 | 5524.30 | |
| Depth to ground-water, ft | 24.9 | 24.85 | 24.32 | 24.02 | |
| Elevation of groundwater, ft | 5499.31 | 5499.45 | 5499.98 | 5500.28 | |

*Summary includes only pollutants determined to be present at concentrations greater than detection limits.

**Laboratory technique error detected.

low

TABLE 4.5
RCRA 3013
GROUNDWATER RESULTS SUMMARY*
MW-5

| Parameter | 3/26/86 | 6/23/86 | 9/18/86 | 12/16/86 | Nominal Detection Limits |
|-----------------------------------|---------|---------|---------|----------|--------------------------|
| Laboratory Analysis (mg/l) | | | | | |
| Cyanide | | 0.2 | 0.24 | | 0.01 |
| Total phenols | 0.006 | 0.007 | 0.034 | 0.021 | 0.001 |
| TOC | 14 | 21 | 20 | 9 | 0.1 |
| TDS | 3,840 | 3,778 | 3,184 | 3,788 | 1 |
| Chloride | 1,100** | 1,339.6 | 1,151 | 1,118 | 1 |
| Sulfate | 14** | 1,800 | 1,237 | 1,132 | 1 |
| Antimony | | | | 0.5 | 0.01 |
| Arsenic | | 0.087 | 0.07 | | 0.05 |
| Cadmium | 0.10 | | | | 0.01 |
| Lead | 0.16 | 0.055 | | | 0.05 |
| Nickel | 0.10 | | 0.09 | 0.07 | 0.06 |
| Selenium | | 0.071 | 0.03 | 0.03 | 0.01 |
| Zinc | 0.012 | 0.02 | 0.02 | 0.016 | 0.01 |
| Barium | | | | 0.01 | 0.005 |
| Manganese | | 0.025 | | | 0.005 |
| Aluminum | | 2.75 | | 4.34 | 0.05 |
| Boron | | | | 0.24 | 0.01 |
| Molybdenum | | | | 0.08 | 0.01 |
| Fluoride | | 0.3 | | 0.580 | 0.01 |
| Nitrate-nitrogen | | 12.5 | | 36 | 0.1 |
| Physical Measurements | | | | | |
| pH, field, s.u. | 7.23 | 7.18 | 7.19 | 7.28 | |
| Conductivity | | 5400 | 6000 | 5700 | |
| T.O.C. elevation, ft | 5545.10 | 5545.10 | 5545.10 | 5545.10 | |
| Depth to groundwater, ft | 40.86 | 40.97 | 41.58 | 44.69 | |

TABLE 4.5 (Continued)

RCRA 3013
GROUNDWATER RESULTS SUMMARY*
MW-5

| Parameter | 3/26/86 | 6/23/86 | 9/18/86 | 12/16/86 | Nominal Detection Limits |
|------------------------------|---------|---------|---------|----------|--------------------------|
| Elevation of groundwater, ft | 5404.24 | 5504.13 | 5503.52 | 5500.41 | |

*Summary includes only pollutants determined to be present at concentrations greater than detection limits.

**Laboratory technique error detected.

TABLE 4.6
RCRA 3013
GROUNDWATER RESULTS SUMMARY*
MW-7

| Parameter | 3/26/86 | 6/23/86 | 9/18/86 | 12/16/86 | Nominal Detection Limits |
|-----------------------------------|---------|---------|---------|----------|--------------------------|
| Laboratory Analysis (mg/l) | | | | | |
| Cyanide | | 0.25 | 0.10 | | 0.01 |
| Total phenols | | 0.006 | 0.036 | 0.025 | 0.001 |
| TOC | 11 | 4 | 4 | 2 | 0.1 |
| TDS | 6,076 | 6,406 | 6,348 | 6,940 | 1 |
| Chloride | 30** | 79.9 | 20 | 29 | 1 |
| Sulfate | 5.5** | 2,400 | 5,802 | 3,630 | 1 |
| Benzene | 0.015 | | 0.058 | 0.009 | 0.001 |
| Toluene | 0.053 | | 0.006 | | 0.001 |
| Ethylbenzene | 0.007 | | 0.004 | | 0.001 |
| Antimony | | | | 0.83 | 0.01 |
| Arsenic | | 0.36 | 0.22 | | 0.05 |
| Cadmium | 0.050 | 0.030 | | 0.02 | 0.01 |
| Chromium | | 0.052 | | 0.08 | 0.05 |
| Lead | | 0.24 | 0.05 | 0.26 | 0.05 |
| Nickel | 0.08 | 0.07 | 0.08 | 0.07 | 0.06 |
| Selenium | | 0.65 | 0.36 | 0.09 | 0.01 |
| Silver | | 0.060 | | | 0.05 |
| Zinc | 0.018 | 0.016 | 0.02 | 0.017 | 0.001 |
| 4,6-Dinitro-o-cresol | 0.013 | | | | 0.001 |
| Benzo(a)anthracene | | 0.001 | | | 0.001 |
| Chrysene | | | | 0.002 | 0.001 |
| 4-Nitrophenol | | | 0.007 | | 0.001 |
| P-chloro-m-cresol | | | 0.001 | | 0.001 |
| Benzo(k)fluoranthene | | | | 0.001 | 0.001 |

TABLE 4.6 (Continued)

RCRA 3013
GROUNDWATER RESULTS SUMMARY*
MW-7

| Parameter | 3/26/86 | 6/23/86 | 9/18/86 | 12/16/86 | Nominal Detection Limits |
|------------------------------|---------|---------|---------|----------|--------------------------|
| Physical Measurements | | | | | |
| pH, field, s.u. | 12.15 | 11.08 | 11.35 | 10.58 | |
| Conductivity | | 8100 | 8600 | 8000 | |
| T.O.C. elevation, ft | 5524.09 | 5524.09 | 5524.09 | 5524.09 | |
| Depth to groundwater, ft | 26.07 | 51.00 | 31.30 | 24.68 | |
| Elevation of groundwater, ft | 5498.02 | 5473.09 | 5492.79 | 5499.41 | |

*Summary includes only pollutants determined to be present at concentrations greater than detection limits.

**Laboratory technique error detected.

TABLE 4.7
RCRA 3013
GROUNDWATER RESULTS SUMMARY*
MW-8

| Parameter | 3/26/86 | 6/23/86 | 9/18/86 | 12/16/86 | Nominal Detection Limits |
|-----------------------------------|---------|---------|---------|----------|--------------------------|
| Laboratory Analysis (mg/l) | | | | | |
| Cyanide | | | | 0.1 | 0.01 |
| Total phenols | | 0.005 | 0.097 | 0.042 | 0.001 |
| TOC | 5 | 13 | 8 | 8 | 0.1 |
| TDS | 806 | 2,910 | 2,284 | 3,450 | 1 |
| Chloride | 160** | 839.7 | 576 | 913 | 1 |
| Sulfate | 4.0** | 1,500 | 586 | 1,270 | 1 |
| Ethylbenzene | 0.107 | | | | 0.01 |
| Antimony | | | | 0.67 | 0.01 |
| Arsenic | | 0.072 | 0.03 | | 0.05 |
| Lead | | 0.055 | | | 0.05 |
| Nickel | | 0.86 | 0.21 | 0.43 | 0.06 |
| Selenium | | 0.21 | | 0.04 | 0.01 |
| Zinc | | 0.020 | 0.02 | 0.016 | 0.01 |
| 4-Nitrophenol | | | 0.008 | | 0.01 |
| Physical Measurements | | | | | |
| pH, field, s.u. | 7.86 | 7.26 | 7.47 | 7.44 | |
| Conductivity | | 4400 | 4000 | 5000 | |
| T.O.C. elevation, ft | 5531.12 | 5531.12 | 5531.12 | 5531.12 | |
| Depth to groundwater, ft | 29.15 | 29.08 | 29.00 | 29.02 | |
| Elevation of groundwater, ft | 5501.97 | 5502.04 | 5502.12 | 5502.1 | |

*Summary includes only pollutants determined to be present at concentrations greater than detection limits.

**Laboratory technique error detected.

TABLE 4.8
 RCRA 3013
 GROUNDWATER RESULTS SUMMARY*
 MW-9

| Parameter | 3/26/86 | 6/23/86 | 9/18/86 | 12/16/86 | Nominal Detection Limits |
|-----------------------------------|---------|---------|---------|----------|--------------------------|
| Laboratory Analysis (mg/l) | | | | | |
| Cyanide | | 0.4 | | | 0.01 |
| Total phenols | 0.304 | 0.372 | 0.17 | 0.16 | 0.001 |
| TOC | 143 | 1,809 | 240 | 275 | 0.1 |
| TDS | 2,360 | 1,718 | 1,428 | 1,684 | 1 |
| Chloride | 149** | 1,009.7 | 89 | 109 | 1 |
| Sulfate | 13.0** | 114 | | 20 | 1 |
| Benzene | 7.4 | 4 | 17.7 | 1.49 | 0.001 |
| Toluene | 6.3 | 1.7 | 10.6 | 0.754 | 0.001 |
| Ethylbenzene | 3.2 | 0.71 | 0.015 | 0.504 | 0.001 |
| Antimony | | | | 0.4 | 0.01 |
| Arsenic | | | 0.02 | | 0.05 |
| Lead | | 0.059 | | | 0.05 |
| Nickel | 0.30 | 0.25 | 0.13 | 0.16 | 0.06 |
| Selenium | | 0.040 | | 0.03 | 0.01 |
| Zinc | 0.012 | 0.015 | 0.05 | 0.011 | 0.01 |
| 2,4-Dimethyl-phenol | 0.160 | 0.150 | | | 0.001 |
| Phenol | 0.149 | 0.170 | 0.013 | 0.133 | 0.001 |
| Fluorene | 0.012 | | | | 0.001 |
| 4-Nitrophenol | | | 1.10 | | 0.001 |
| Acenaphthalene | | | 0.028 | | 0.001 |
| Benzo(a)anthracene | | | 0.007 | | 0.001 |
| Pyrene | | | 0.010 | | 0.001 |
| Naphthalene | | | | 0.029 | 0.001 |

TABLE 4.8 (Continued)

RCRA 3013
GROUNDWATER RESULTS SUMMARY*
MW-9

| Parameter | 3/26/86 | 6/23/86 | 9/18/86 | 12/16/86 | Nominal Detection Limits |
|------------------------------|---------|---------|---------|----------|--------------------------|
| Physical Measurements | | | | | |
| pH, field, s.u. | 7.01 | 6.98 | 6.89 | 6.91 | |
| Conductivity | | 2500 | 2200 | 2600 | |
| T.O.C. elevation, ft | 5519.70 | 5519.70 | 5519.70 | 5519.70 | |
| Depth to groundwater, ft | 21.50 | 20.23 | 20.13 | 20.55 | |
| Elevation of groundwater, ft | 5498.20 | 5499.47 | 5499.57 | 5499.15 | |

*Summary includes only pollutants determined to be present at concentrations greater than detection limits.

**Laboratory technique error detected.

TABLE 4.9
RCRA 3013
GROUNDWATER RESULTS SUMMARY*
MW-10

| Parameter | 3/26/86 | 6/23/86 | 9/18/86 | 12/16/86 | Nominal Detection Limits |
|-----------------------------------|---------|---------|---------|----------|--------------------------|
| Laboratory Analysis (mg/l) | | | | | |
| Cyanide | | | 0.050 | | 0.01 |
| Total phenols | 0.147 | 0.186 | 0.065 | 0.055 | 0.001 |
| TOC | 34 | 76 | 125 | 114 | 0.1 |
| TDS | 1,546 | 2,820 | 2,408 | 3,272 | 1 |
| Chloride | 245** | 569.8 | 587 | 457 | 1 |
| Sulfate | 5.3** | 165 | | 10 | 1 |
| Benzene | 0.093 | | 0.041 | 14.1 | 0.002 |
| Toluene | | | 0.054 | 7.4 | 0.001 |
| Ethylbenzene | | | | 0.03 | 0.001 |
| Antimony | | | | 0.56 | 0.01 |
| Arsenic | | 0.053 | | | 0.05 |
| Beryllium | | | | 0.04 | 0.01 |
| Cadmium | 0.020 | | | | 0.01 |
| Lead | | 0.059 | 0.05 | | 0.05 |
| Nickel | 0.08 | 0.25 | 0.18 | 0.08 | 0.06 |
| Selenium | | 0.040 | 0.071 | 0.03 | 0.01 |
| Zinc | | 0.015 | 0.16 | 0.01 | 0.01 |
| 2,4-Dimethyl-phenol | 0.025 | | | | 0.001 |
| 4,6-Dinitro-o-cresol | 0.020 | | | | 0.001 |
| Phenol | 0.090 | | | | 0.001 |
| Anthracene | 0.039 | | | | 0.001 |
| Fluoranthene | 0.034 | | | | 0.001 |
| Fluorene | 0.033 | | | | 0.001 |
| Pyrene | 0.030 | | | | 0.001 |
| 2-Nitrophenol | | | 0.002 | | 0.001 |
| 4-Nitrophenol | | | 0.016 | | 0.001 |
| Naphthalene | | | | 0.004 | 0.001 |

TABLE 4.9 (Continued)

RCRA 3013
GROUNDWATER RESULTS SUMMARY*
MW-10

| Parameter | 3/26/86 | 6/23/86 | 9/18/86 | 12/16/86 | Nominal Detection Limits |
|------------------------------|---------|---------|---------|----------|--------------------------|
| Physical Measurements | | | | | |
| pH, field, s.u. | 7.07 | 7.08 | 6.93 | 7.05 | |
| Conductivity | | 4400 | 4800 | 5100 | |
| T.O.C. elevation, ft | 5516.86 | 5516.86 | 5516.86 | 5516.86 | |
| Depth to groundwater, ft | 19.20 | 18.75 | 18.11 | 17.56 | |
| Elevation of groundwater, ft | 5497.66 | 5498.11 | 5498.75 | 5499.3 | |

*Summary includes only pollutants determined to be present at concentrations greater than detection limits.

**Laboratory technique error detected.

of these compounds is believed to suggest the possibility of laboratory error - particularly when the reported concentration of the substance is very near detection limits.

disagree

A good example of the singularity of results is seen in Table 4.6, which reports a summary of monitoring well 7 results. Every hydrocarbon substance detected was at low levels close to the detection limit. Laboratory error may be responsible for detection of these substances.

The monitoring well sampling method is also important in assessing the significance of sample results. Basically, the technique involved BRC staff collection of a sample of water such that the air/liquid interface was acquired in the sample. This process maximized the possibility of acquiring any free-phase hydrocarbons and thus the highest concentrations in the sample. Therefore, sample concentrations for each sampling effort constitute a worst-case assessment of the well water when floating separate-phase contaminants are present.

The significance of concentrations measured at the BRC monitoring wells should be judged from two different standpoints: (1) water use, and (2) subsurface hydrocarbon contamination. From a water use standpoint, the concentrations of pollutants under the BRC site are insignificant. Groundwater is not currently used, and there are no plans for its use in the future.

The monitoring well sampling results include an indicator parameter in total organic carbon (TOC). This parameter can be useful in identifying suspected locations of significant concentrations of subsurface hydrocarbon since TOC values are generally higher in samples having free product.

A review of Tables 4.1 through 4.10 show that TOC values measured in monitoring well 4 were higher than all wells except 9 and 10. Well 9 has exhibited an indication of films of hydrocarbon material at the surface (reported by BRC staff on January 27, 1987). The fact that the TOC results from this well are the highest is consistent with an indication of films in this well. Films have also been indicated in monitoring well 10 during the third and fourth quarter sampling efforts recorded by

BRC staff January 27, 1987. This evidence is consistent with the highest TOC values recorded for these two sampling efforts. The proximity of well 10 to well 4 and the similarity of their respective TOC values during periods of indications of floating hydrocarbon evidence suggest they may be influenced by the same source of hydrocarbon material.

CHAPTER 5

GROUNDWATER MONITORING WELL WATER LEVEL MEASUREMENTS

BACKGROUND

In accordance with the EPA-approved work plan, water levels have been measured by Bloomfield staff monthly at each of the groundwater monitoring wells located on the Bloomfield refinery property. Water level measurements commenced in January 1986. Since installation of monitoring wells 7 through 10 was completed in March 1986, water level measurements for these wells commenced in March.

In addition to water level measurements made in the wells, measurements were also made on Hammond Ditch at its intersection with Sullivan Road and at a walkway bridge on refinery property. These water levels and the well water level measurements are presented in Figures 5.1 through 5.16. Hammond Ditch measurement locations are also indicated on the figures.

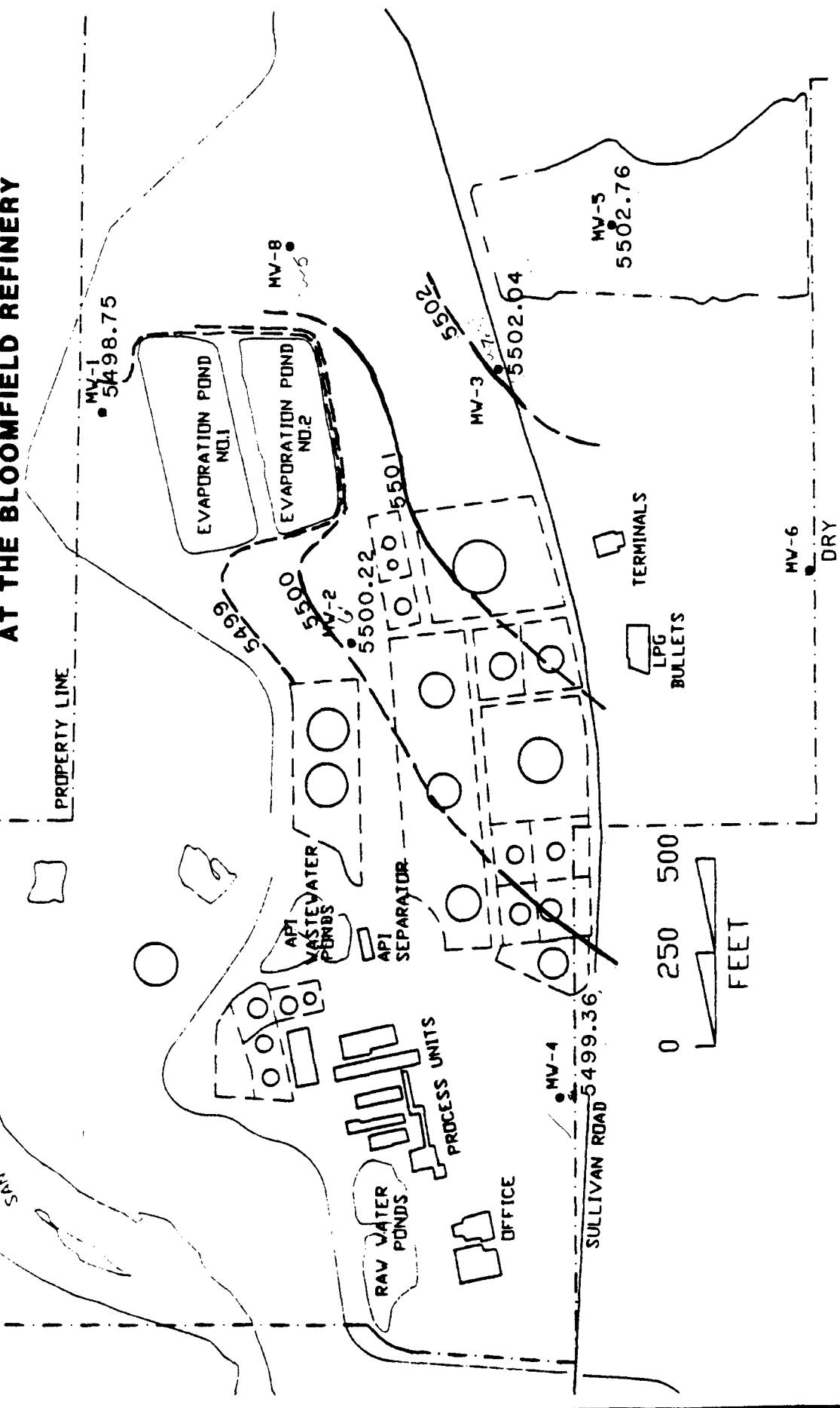
METHODOLOGY

Two approaches were employed to evaluate the monthly water level measurements. In the first approach, groundwater elevations measured at each well were plotted on a map of the BRC site. Then groundwater contours were drawn on these maps based on the relative orientation of measured elevations. The water levels shown by monitoring well 7 were ignored in this process because they represented water levels in the Nacimiento rather than the cobble zone (represented by the other wells). The contours were drawn using both solid and dashed lines. Solid lines constitute an interpolative estimation of groundwater gradient for the given data of groundwater elevations. The dashed lines constitute a presumption of the influences of Hammond Ditch, the evaporation pond, and spray irrigation fields on the groundwater regime. These lines also assume that no irregularities in flow patterns exist in the subsurface beneath the process and refined product storage areas. These maps are presented as Figures 5.1 through 5.16.

NOF
#4200
#70
area

FIGURE 5.1
**WATER LEVEL MEASUREMENTS WITH
 ESTIMATED GROUNDWATER CONTOURS FOR
 AT THE BLOOMFIELD REFINERY**

1/24/86



**FIGURE 5.2
WATER LEVEL MEASUREMENTS WITH
ESTIMATED GROUNDWATER CONTOURS FOR
2/20/86**

AT THE BLOOMFIELD REFINERY

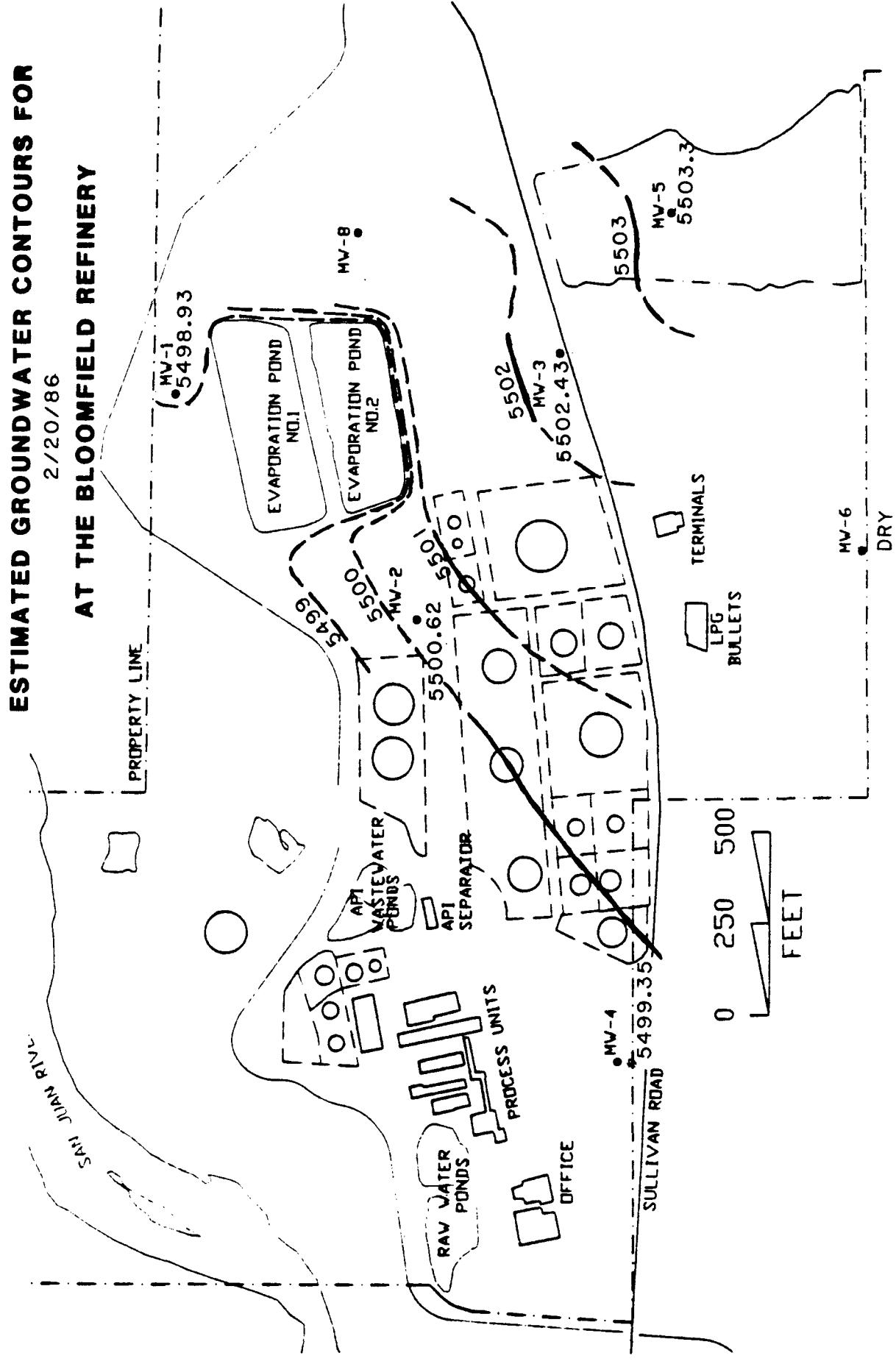


FIGURE 5.3
**WATER LEVEL MEASUREMENTS WITH
 ESTIMATED GROUNDWATER CONTOURS FOR
 AT THE BLOOMFIELD REFINERY**

3/21/86

PROPERTY LINE

SULLIVAN RIVER

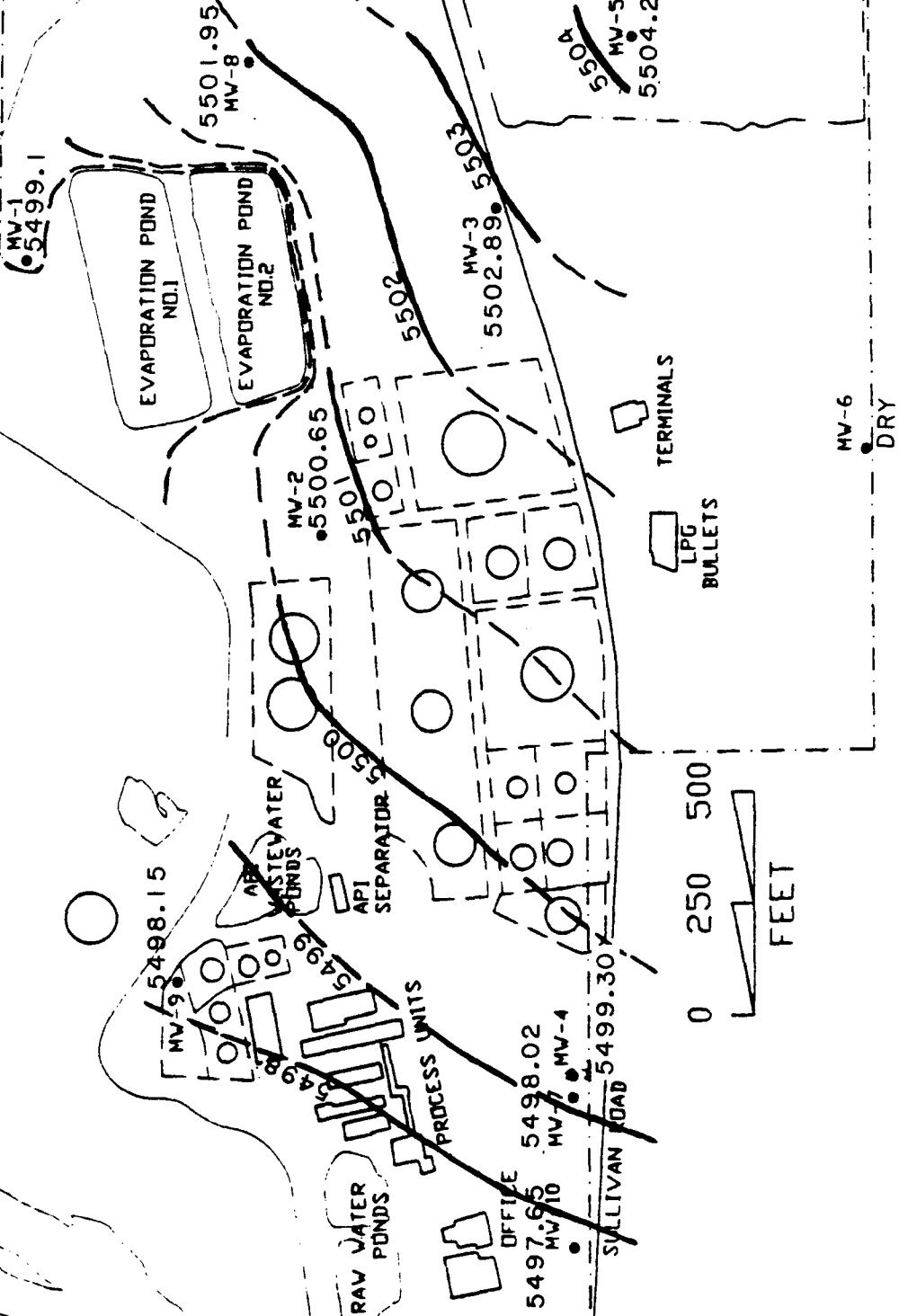


FIGURE 5.4
**WATER LEVEL MEASUREMENTS WITH
 ESTIMATED GROUNDWATER CONTOURS FOR
 AT THE BLOOMFIELD REFINERY**
 3/26/86

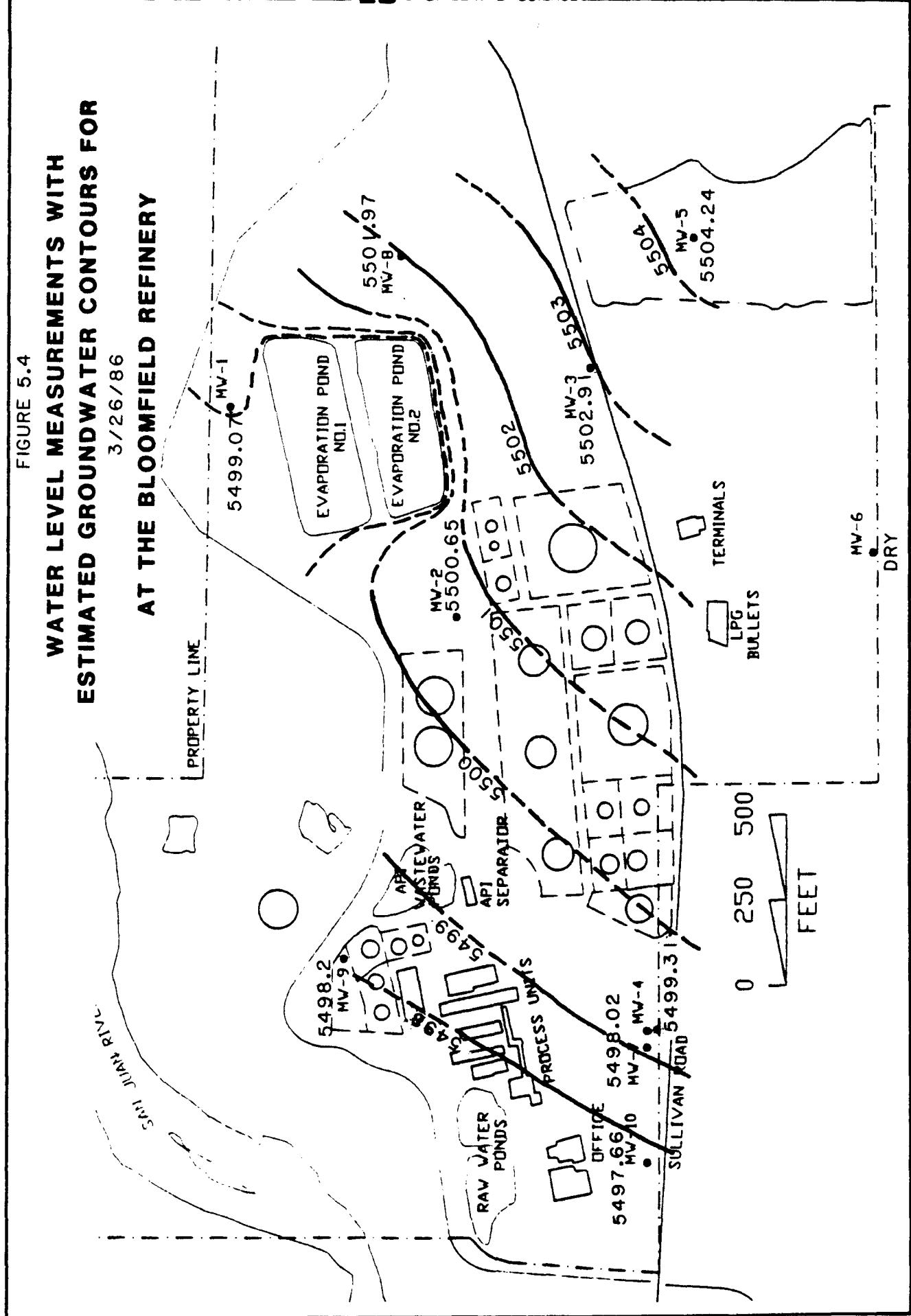


FIGURE 5.5
**WATER LEVEL MEASUREMENTS WITH
 ESTIMATED GROUNDWATER CONTOURS FOR
 AT THE BLOOMFIELD REFINERY**

4/4/86

PROPERTY LINE

SALT RIVER

MW-1 99.07

EVAPORATION POND NO.1

EVAPORATION POND NO.2

5501.86 MW-8

5500.52 MW-2

5502.98 MW-3

5504.57 MW-5

5498.22 MW-9

5498.24 API SEPARATOR

5497.6 MW-10

5498.77 MW-7

5499.2 MW-4

5496.5 MW-6

SULLIVAN ROAD

OFFICE

RAW WATER PONDS

API SEPARATORS

PROCESS UNITS

TERMINALS

LPG BULLETS

DRY

SURFACE WATER SAMPLING SITE

FEET

MW-6

DRY

FIGURE 5.6
WATER LEVEL MEASUREMENTS WITH
ESTIMATED GROUNDWATER CONTOURS FOR
4/18/86
AT THE BLOOMFIELD REFINERY

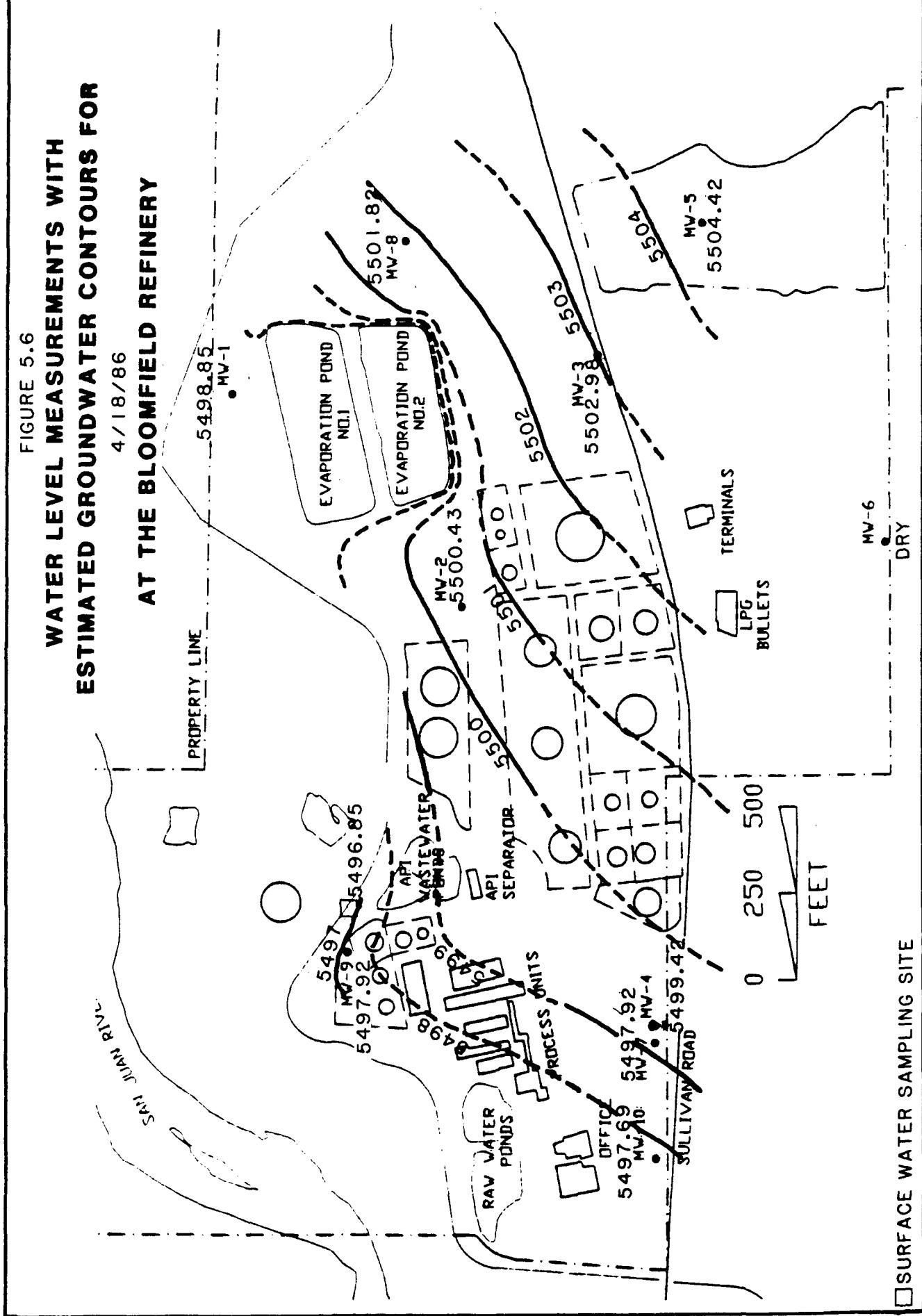


FIGURE 5.7
**WATER LEVEL MEASUREMENTS WITH
 ESTIMATED GROUNDWATER CONTOURS FOR
 5/5/86**
AT THE BLOOMFIELD REFINERY

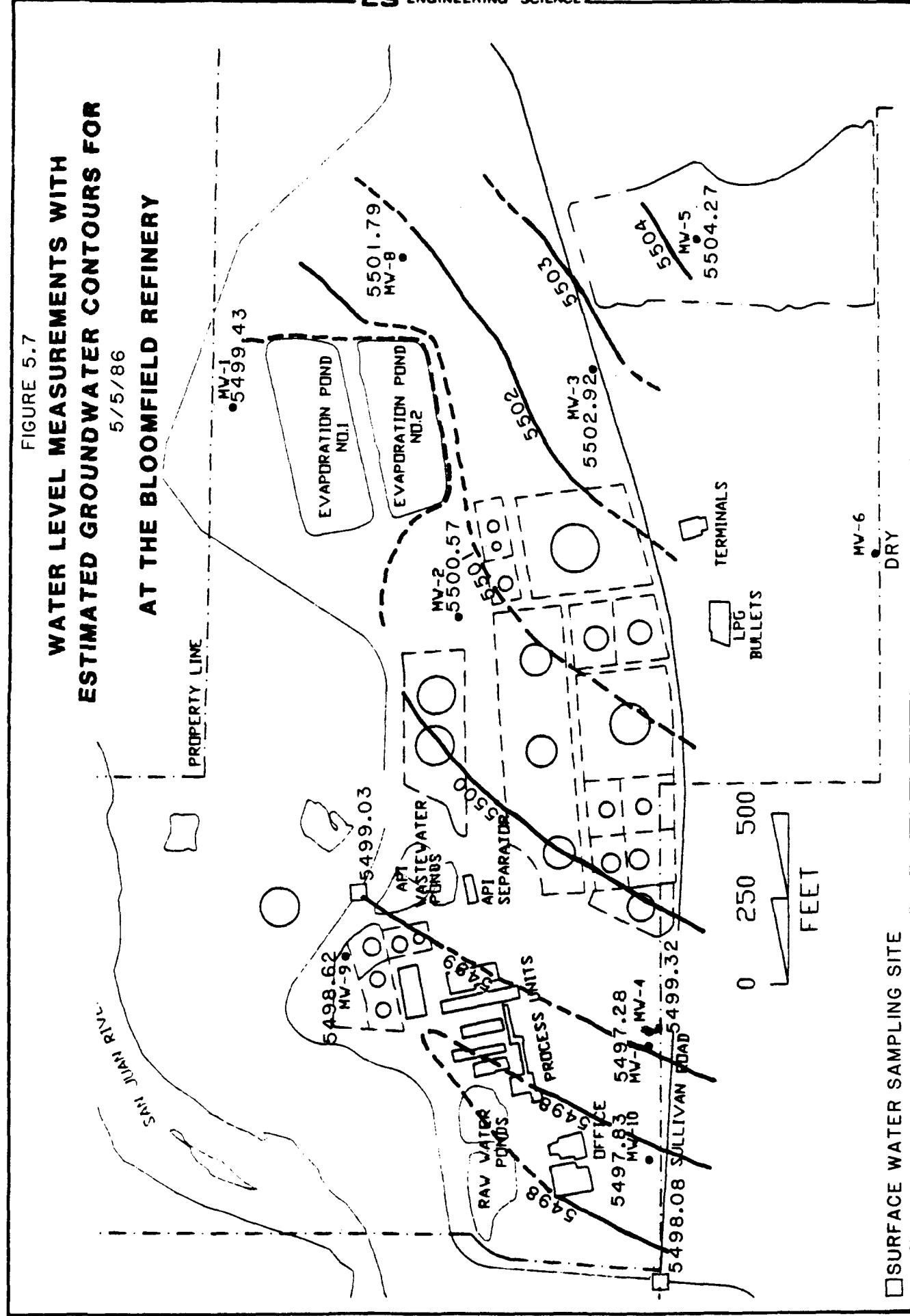
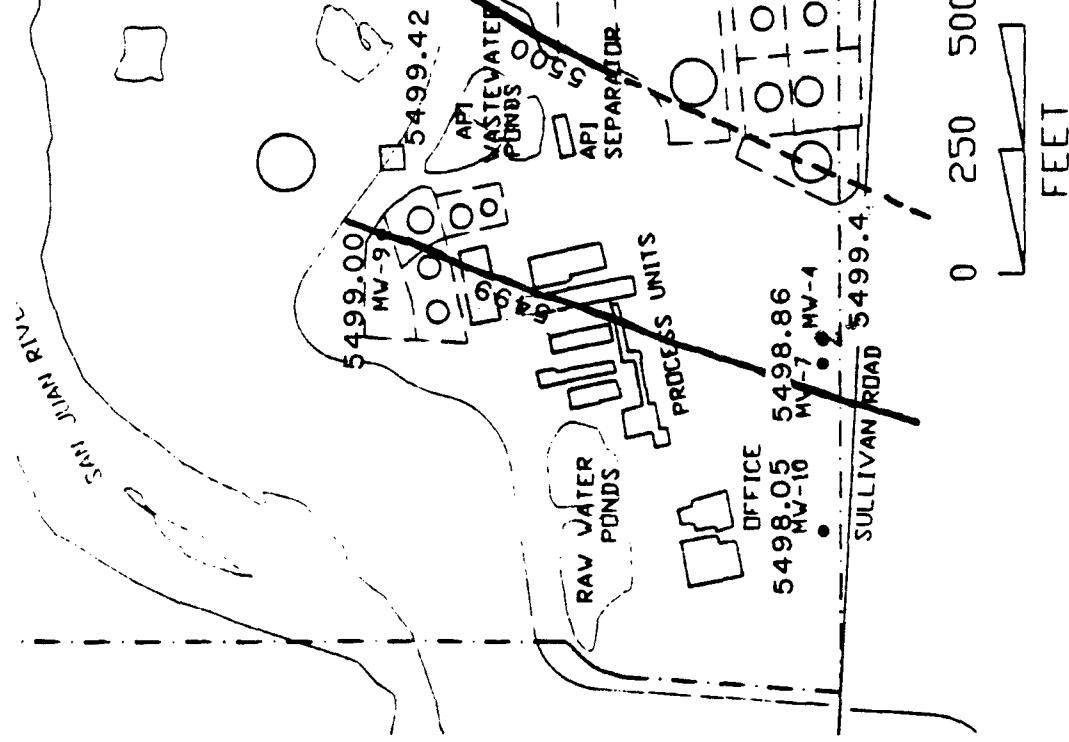


FIGURE 5.8
WATER LEVEL MEASUREMENTS WITH
ESTIMATED GROUNDWATER CONTOURS FOR
AT THE BLOOMFIELD REFINERY

5/21/86

MW-1
5500.05MW-8
5501.83EVAPORATION POND
NO.1EVAPORATION POND
NO.2MW-2
5500.82MW-3
5502.85MW-2
5502MW-3
5503MW-4
5504
MW-5
5504.35MW-6
DRYTERMINALS
LPG
BULLETS0
250
500
FEET

□ SURFACE WATER SAMPLING SITE

FIGURE 5.9
WATER LEVEL MEASUREMENTS WITH
ESTIMATED GROUNDWATER CONTOURS FOR
AT THE BLOOMFIELD REFINERY

6/4/86

MW-1

MW-1

5500.41

EVAPORATION POND
NO.1EVAPORATION POND
NO.2

MW-2

5500.93

MW-2

5

FIGURE 5.10
**WATER LEVEL MEASUREMENTS WITH
 ESTIMATED GROUNDWATER CONTOURS FOR
 AT THE BLOOMFIELD REFINERY**

7/8/86

PROPERTY LINE

SILVER RIVER

WASTED WATER PONDS

SEPARATOR

PROCESS UNITS

OFFICE

MV-9

MV-10

MV-7

MV-4

MV-6

MV-3

MV-2

MV-1

MV-8

MV-5

MV-2.2

MV-03

MV-02.96

MV-03

MV-02

MV-01.34

MV-02.27

MV-03

MV-03.87

MV-06

DRY

SURFACE WATER SAMPLING SITE

LPG

BULLETS

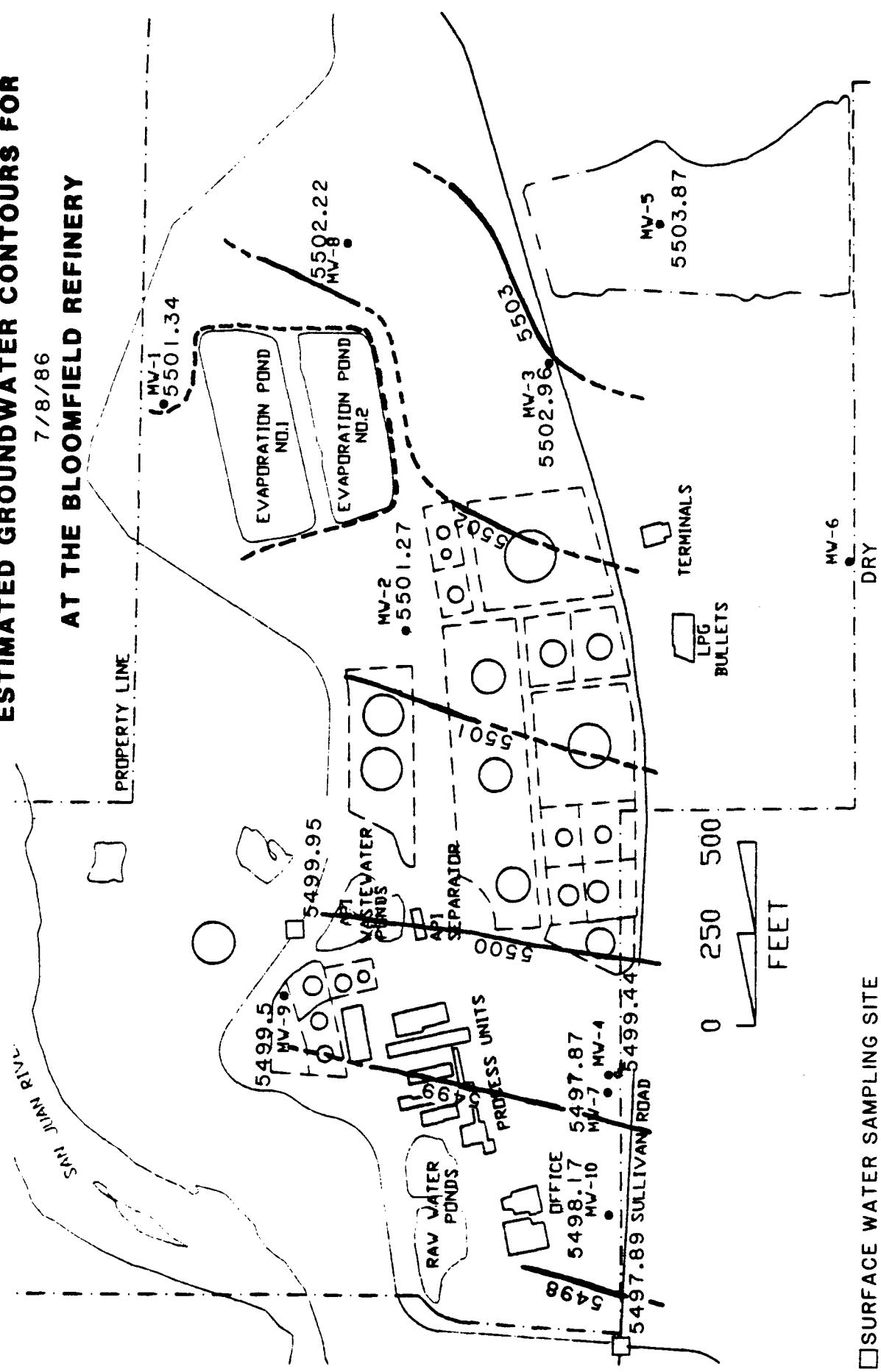
TERMINALS

0

250

500

FEET



**FIGURE 5.11
WATER LEVEL MEASUREMENTS WITH
ESTIMATED GROUNDWATER CONTOURS FOR
AT THE BLOOMFIELD REFINERY**

8/4/86

PROPERTY LINE

MW-1
5500.25MW-8
5502.12MW-3
5502.92MW-5
5503.77MW-6
DRY

EVAPORATION POND NO.1

EVAPORATION POND NO.2

MW-2
5501.13

TERMINALS

LPG
BULLETS

0 250 500 FEET

5499.4
MW-9
5499.67API
VACUUM
PUMPS
SEPARATORPROCESS UNITS
RAW WATER PONDSMW-10
5498.37MW-7
5498.77MW-4
5499.67MW-1
5497.84 SULLIVAN ROAD

□ SURFACE WATER SAMPLING SITE

JUAN RIVER

FIGURE 5.12
WATER LEVEL MEASUREMENTS WITH
ESTIMATED GROUNDWATER CONTOURS FOR
AT THE BLOOMFIELD REFINERY

9/2/86

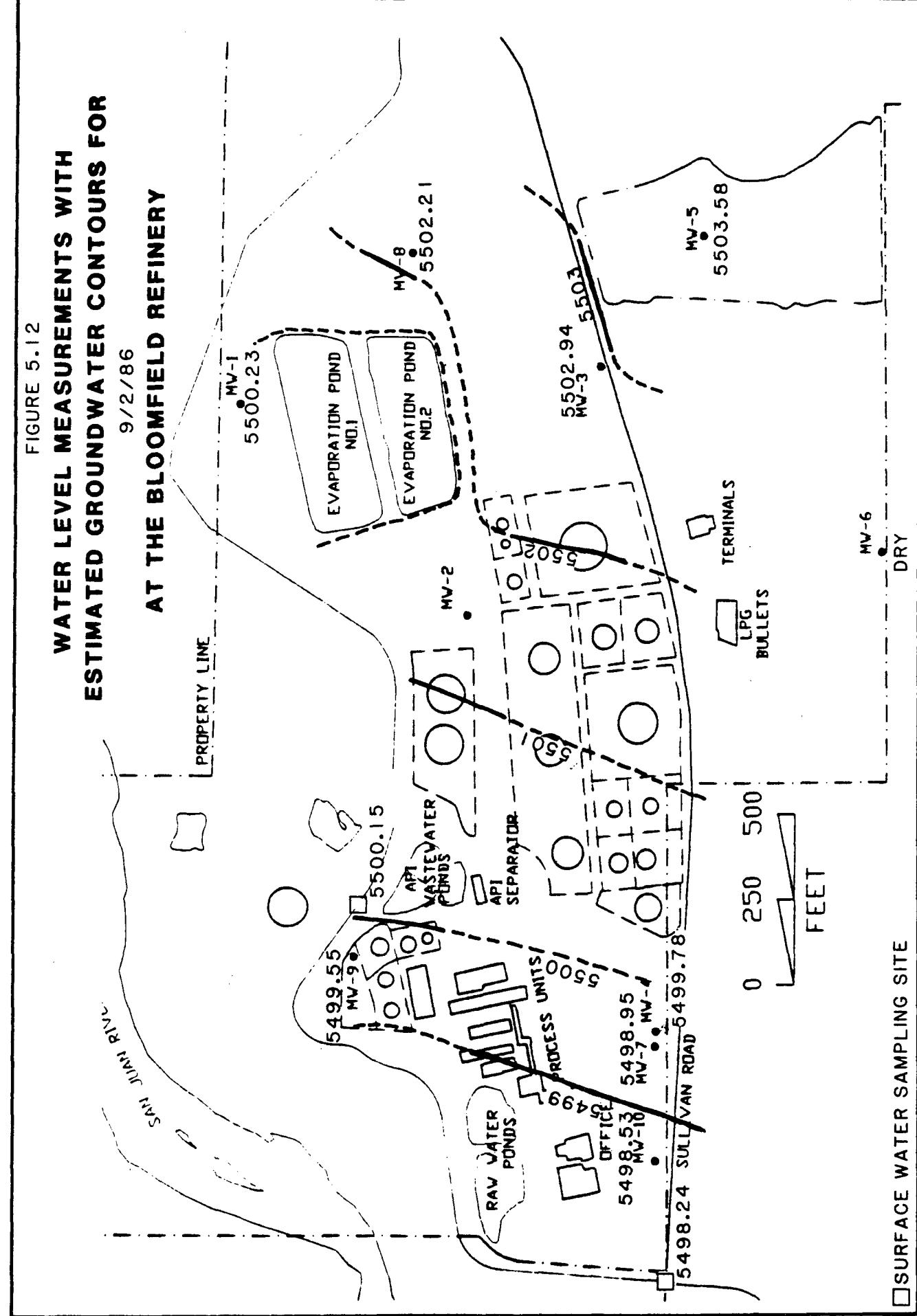


FIGURE 5.13
**WATER LEVEL MEASUREMENTS WITH
 ESTIMATED GROUNDWATER CONTOURS FOR
 10/8/86
 AT THE BLOOMFIELD REFINERY**

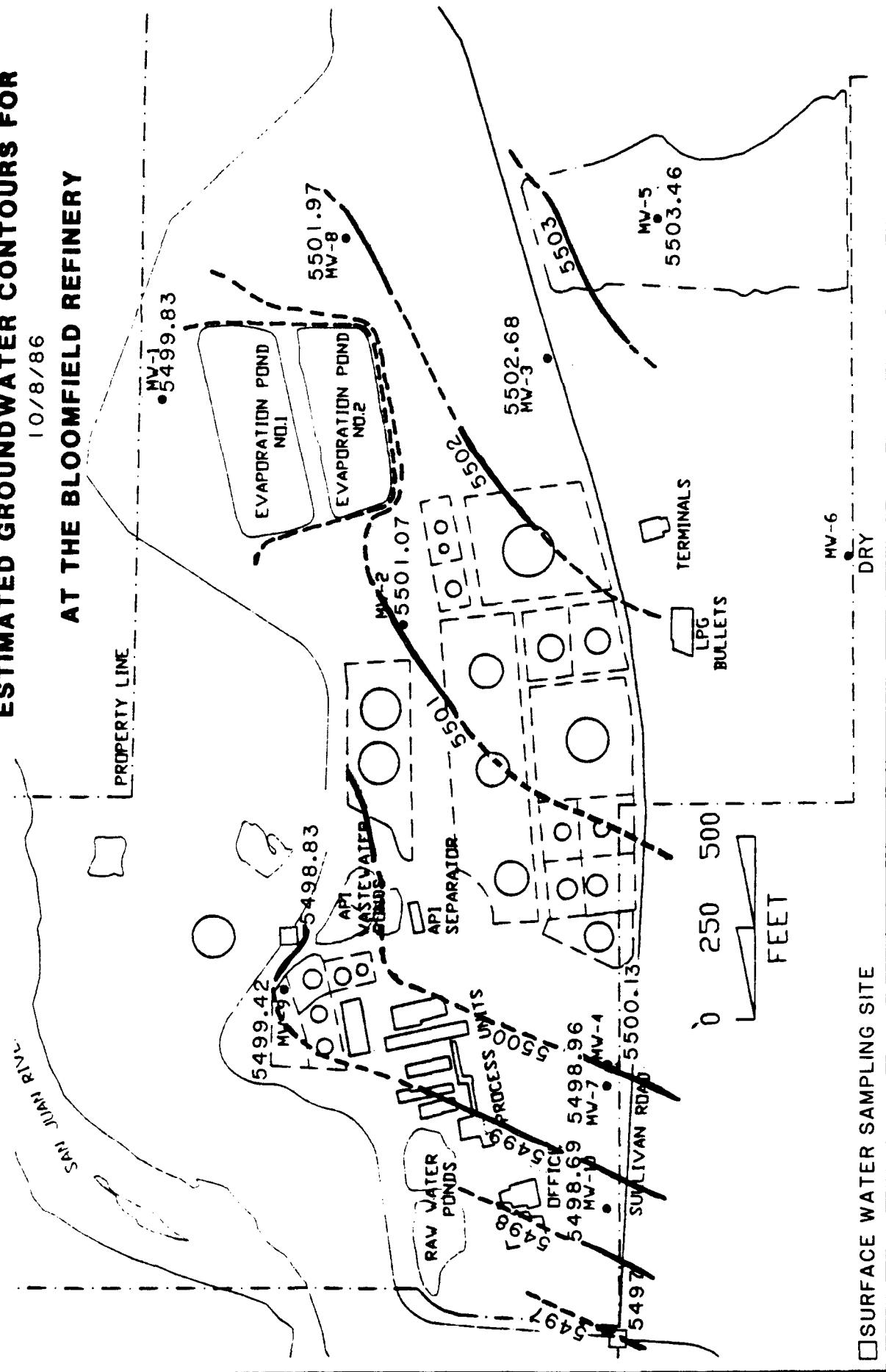
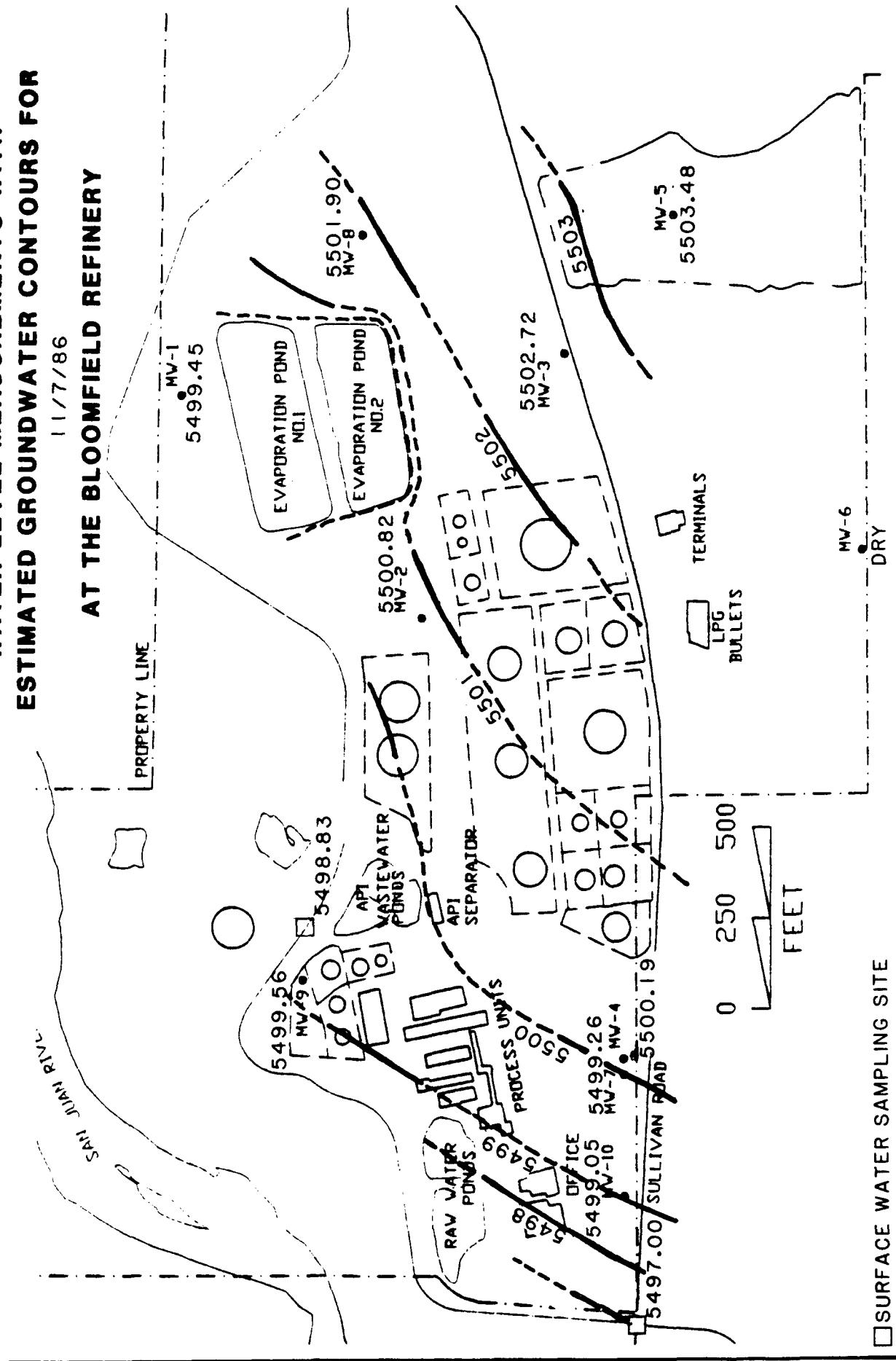


FIGURE 5.14
**WATER LEVEL MEASUREMENTS WITH
 ESTIMATED GROUNDWATER CONTOURS FOR
 AT THE BLOOMFIELD REFINERY**

11/7/86



□ SURFACE WATER SAMPLING SITE

FIGURE 5.15
**WATER LEVEL MEASUREMENTS WITH
 ESTIMATED GROUNDWATER CONTOURS FOR
 AT THE BLOOMFIELD REFINERY**
 12/8/86

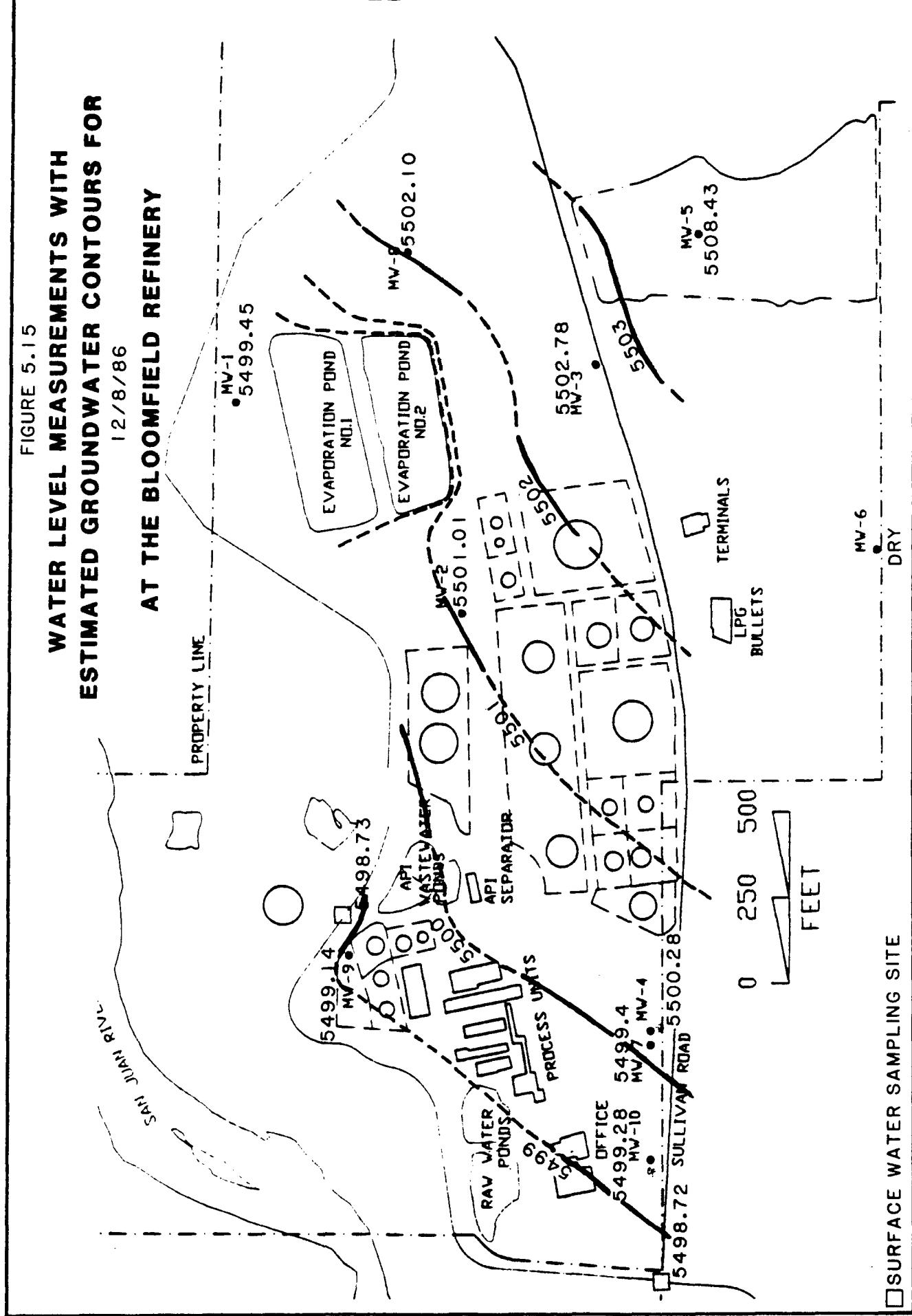
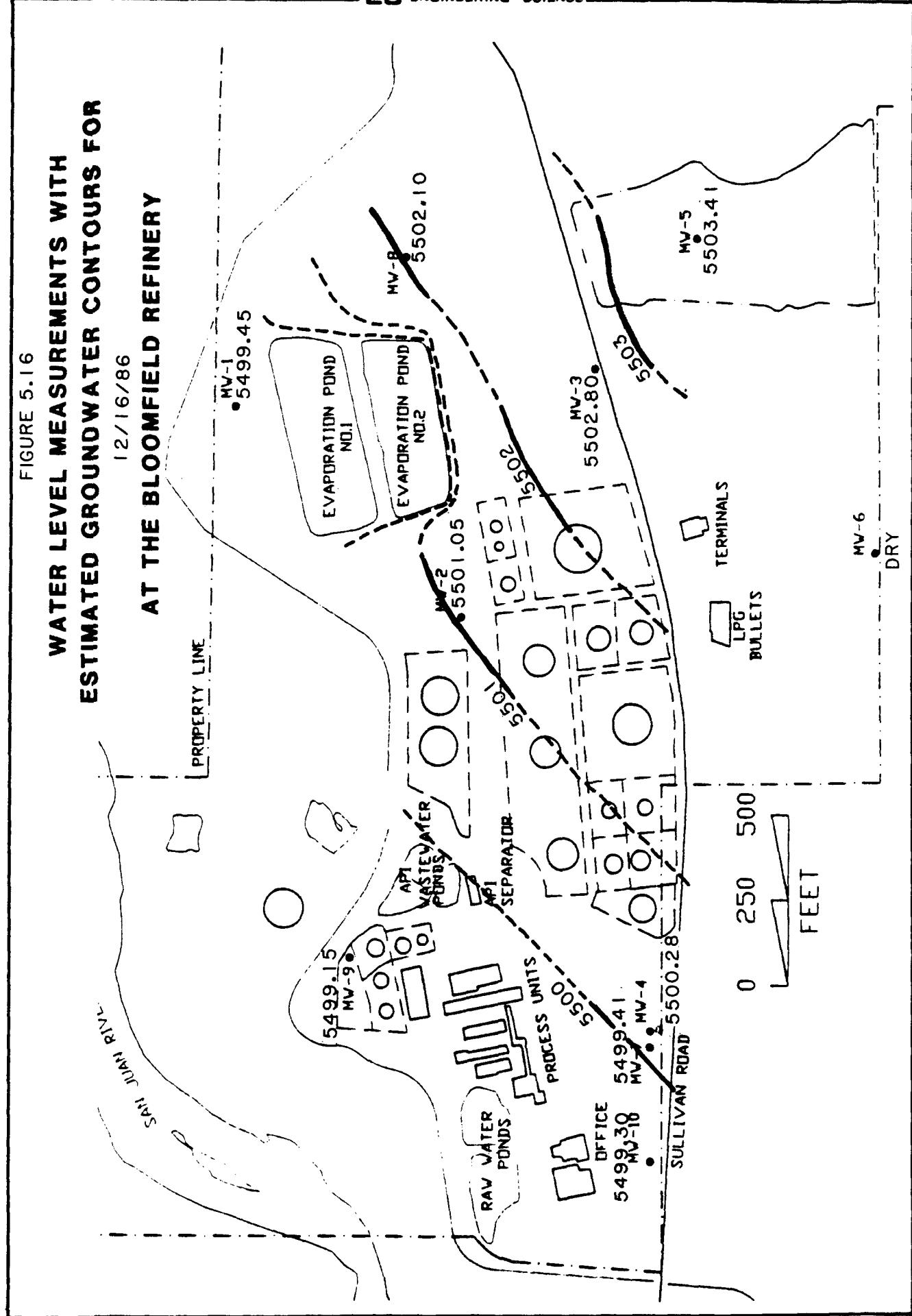


FIGURE 5.16
12/16/86
**WATER LEVEL MEASUREMENTS WITH
ESTIMATED GROUNDWATER CONTOURS FOR
AT THE BLOOMFIELD REFINERY**



In the second approach, time series plots were constructed of groundwater elevations measured in each well. A listing of the plotted elevations is provided in Table 5.1. Time series plots of surface water elevations, determined by BRC staff for Hammond Ditch at both Sullivan Road and a walkway bridge located upstream on BRC property, were superimposed on monitoring wells 1, 4, 9, and 10 due to their proximity to Hammond Ditch. These plots, presented as Figures 5.17 and 5.18, were developed to evaluate the effect of Hammond Ditch on the site's groundwater. Figure 5.19 contains a comparison of the response of monitoring wells 2, 3, 5, 8, and 9. Figure 5.20 compares monitoring wells 4 and 7.

DISCUSSION

Based on the delineation of contours on Figures 5.1 through 5.16, flow through the site occurs generally in a northwesterly direction toward Hammond Ditch. However, there is an insufficient number of wells over the site to obtain more than a generalized picture of flow.

The time series plots presented in Figures 5.17 and 5.18 show that groundwater levels are definitely influenced by flow in Hammond Ditch. In Figure 5.17, water levels in Hammond Ditch at the walkway bridge are shown to be higher than water levels in monitoring well 9. This relationship means that, during irrigation season, groundwater from the site does not flow into Hammond Ditch. Instead, water flows from higher elevations in the refinery to a point where water mounded in the ditch bank precludes further flow toward the ditch. It is believed the mounding effect creates a trough in groundwater contours on the refinery side of the ditch. This trough is believed to be an avenue of transport for groundwater movement downstream along a path parallel to Hammond Ditch. The width of the trough should be locally affected by the area groundwater gradients.

The relationship of the groundwater level in monitoring wells 2, 3, 5, 8, and 9 to seasonal Hammond Ditch flows is presented in Figure 5.19. In this figure, monitoring well 9 is used as a basis of comparison from water levels presented in Figure 5.17. Based on the performance of these

TABLE 5.1

GROUNDWATER ELEVATIONS MEASURED AT BLOOMFIELD REFINING COMPANY

| Date | MW-1 | MW-2 | MW-3 | MW-4 | MW-5 | MW-7 | MW-8 | MW-9 | MW-10 | Hammond Ditch at Sullivan Road | Hammond Ditch at Walkway |
|----------|---------|---------|---------|---------|---------|---------|---------|---------|---------|---|--------------------------------|
| 02/23/85 | 5499.07 | 5500.55 | 5502.15 | 5499.30 | 5502.75 | | | | | | |
| 03/18/85 | 5499.14 | 5500.82 | 5502.55 | 5499.32 | 5503.50 | | | | | | |
| 04/11/85 | 5498.99 | 5500.62 | 5502.73 | 5499.30 | 5503.67 | | | | | | |
| 05/31/85 | 5499.67 | 5500.97 | 5502.74 | 5499.80 | 5503.64 | | | | | | |
| 06/14/85 | 5499.80 | 5500.99 | 5502.83 | 5499.80 | 5503.40 | | | | | | |
| 06/26/85 | 5499.94 | 5500.98 | 5502.49 | 5499.73 | 5503.24 | | | | | | |
| 07/10/85 | 5500.20 | 5500.99 | 5502.48 | 5499.80 | 5503.30 | | | | | | |
| 08/02/85 | 5501.00 | 5501.25 | 5502.48 | 5499.78 | 5503.37 | | | | | | |
| 09/17/85 | 5500.34 | 5501.05 | 5502.25 | 5499.80 | 5503.00 | | | | | | |
| 10/09/85 | 5500.03 | 5500.87 | 5502.42 | 5499.70 | 5503.30 | | | | | | |
| 10/24/85 | 5499.23 | 5500.43 | 5502.28 | 5499.54 | 5503.10 | | | | | | |
| 11/08/85 | 5498.72 | 5500.05 | 5502.20 | 5499.60 | 5503.09 | | | | | | |
| 12/17/85 | 5498.35 | 5499.85 | 5501.85 | 5499.40 | 5502.90 | | | | | | |
| 01/08/86 | 5498.59 | 5500.08 | 5501.85 | 5499.35 | 5502.77 | | | | | | |
| 01/24/86 | 5498.75 | 5500.22 | 5502.04 | 5499.36 | 5502.76 | | | | | | |
| 02/20/86 | 5498.93 | 5500.62 | 5502.43 | 5499.35 | 5503.30 | | | | | | |
| 03/21/86 | 5499.10 | 5500.65 | 5502.89 | 5499.30 | 5504.23 | 5498.02 | 5501.95 | 5498.15 | 5497.65 | | |
| 03/26/86 | 5499.07 | 5500.65 | 5502.91 | 5499.31 | 5504.24 | 5498.02 | 5501.97 | 5498.20 | 5497.66 | | |
| 04/04/86 | 5499.07 | 5500.57 | 5502.98 | 5499.21 | 5504.57 | 5498.77 | 5501.86 | 5498.22 | 5497.60 | 5496.50 | 5498.24 |
| 04/18/86 | 5498.85 | 5500.43 | 5502.98 | 5499.42 | 5504.42 | 5497.92 | 5501.82 | 5498.90 | 5497.69 | | 5496.85 |
| 05/05/86 | 5499.43 | 5500.57 | 5502.92 | 5499.32 | 5504.27 | 5497.28 | 5501.79 | 5498.62 | 5497.83 | 5498.08 | 5499.03 |
| 05/21/86 | 5500.05 | 5500.82 | 5502.85 | 5499.40 | 5504.35 | 5498.86 | 5501.83 | 6599.00 | 5498.05 | 5498.25 | 5499.42 |
| 06/04/86 | 5500.41 | 5500.93 | 5502.95 | 5499.40 | 5504.17 | 5498.85 | 5501.89 | 5499.17 | 5498.15 | 5498.23 | 5499.63 |
| 06/23/86 | 5501.21 | 5501.18 | 5503.05 | 5499.45 | 5504.13 | | | | | | |
| 07/08/86 | 5501.34 | 5501.27 | 5502.96 | 5499.44 | 5503.87 | 5497.87 | 5502.22 | 5499.50 | 5498.17 | 5497.89 | 5499.95 |
| 08/04/86 | 5500.25 | 5501.13 | 5502.92 | 5499.67 | 5503.77 | 5498.77 | 5502.12 | 5499.40 | 5498.37 | 5497.84 | 5499.67 |
| 09/02/86 | 5500.23 | 5501.32 | 5502.94 | 5499.78 | 5503.58 | 5498.95 | 5502.21 | 5499.55 | 5498.53 | 5498.24 | 5500.15 |
| 09/18/86 | 5500.03 | 5501.22 | 5502.77 | 5499.98 | 5503.52 | | | | | | |
| 10/08/86 | 5499.83 | 5501.07 | 5502.68 | 5500.13 | 5503.46 | 5499.96 | 5501.97 | 5499.42 | 5498.69 | 5497.00 | 5498.83 |
| 11/07/86 | 5499.45 | 5500.82 | 5502.72 | 5500.19 | 5503.48 | 5499.26 | 5501.90 | 5499.56 | 5499.05 | 5497.00 | 5498.83 |
| 12/08/86 | 5499.45 | 5501.01 | 5502.78 | 5500.28 | 5503.43 | 5499.40 | 5502.10 | 5499.14 | 5499.28 | 5498.72 | 5498.73 |
| 12/16/86 | 5499.45 | 5501.05 | 5502.80 | 5500.28 | 5503.41 | 5499.41 | 5502.10 | 5499.15 | 5499.30 | | |

NOTES:

MW-6 was dry for the entire time period.

Measurements from project startup through 7/8/86 were not stabilized due to slow recovery from well purging.

FIGURE 5.17
GROUNDWATER ELEVATIONS
BLOOMFIELD REFINING – BLOOMFIELD, NEW MEXICO
MONITORING WELLS 1, 9, HAW, HAS

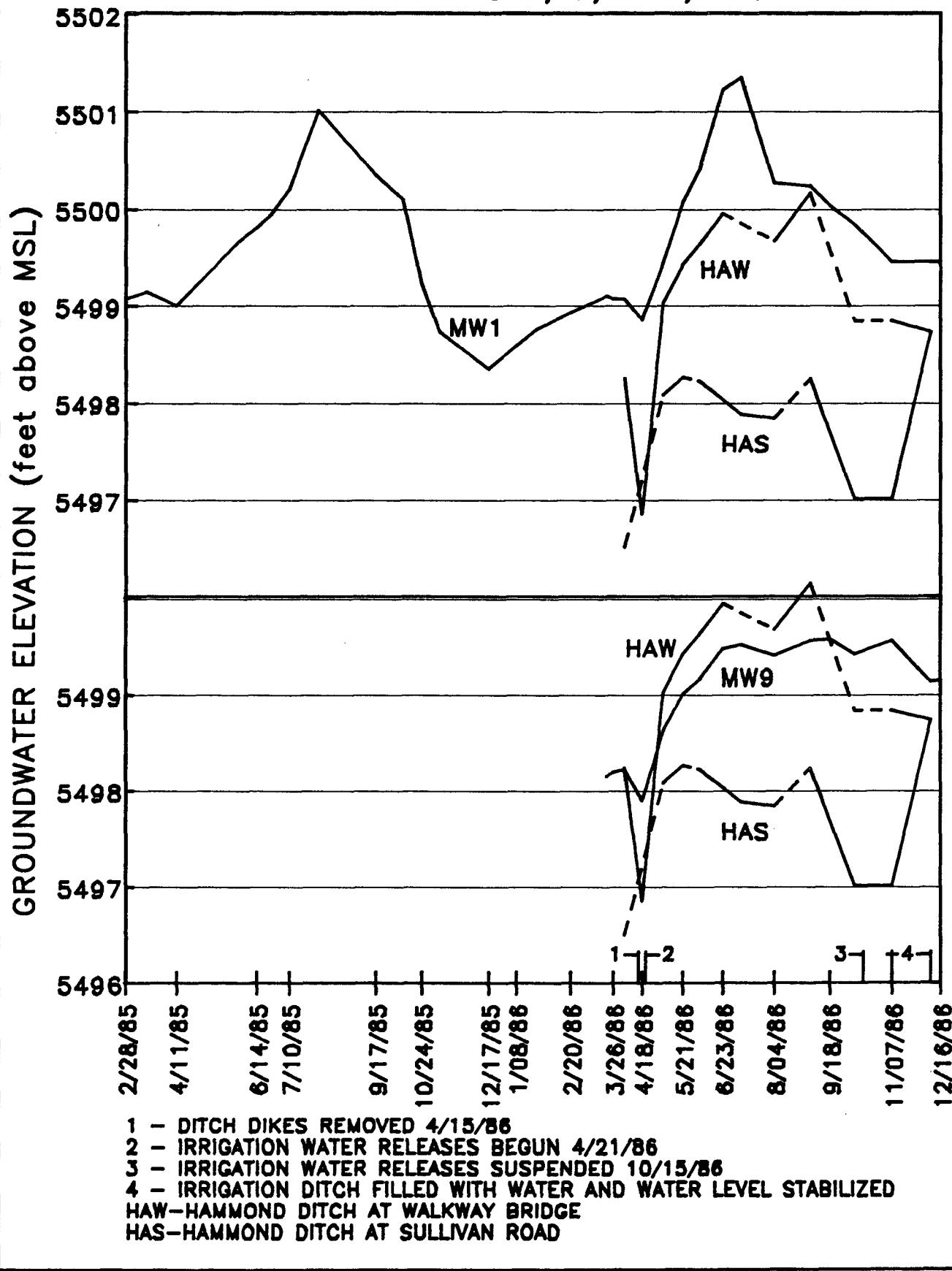
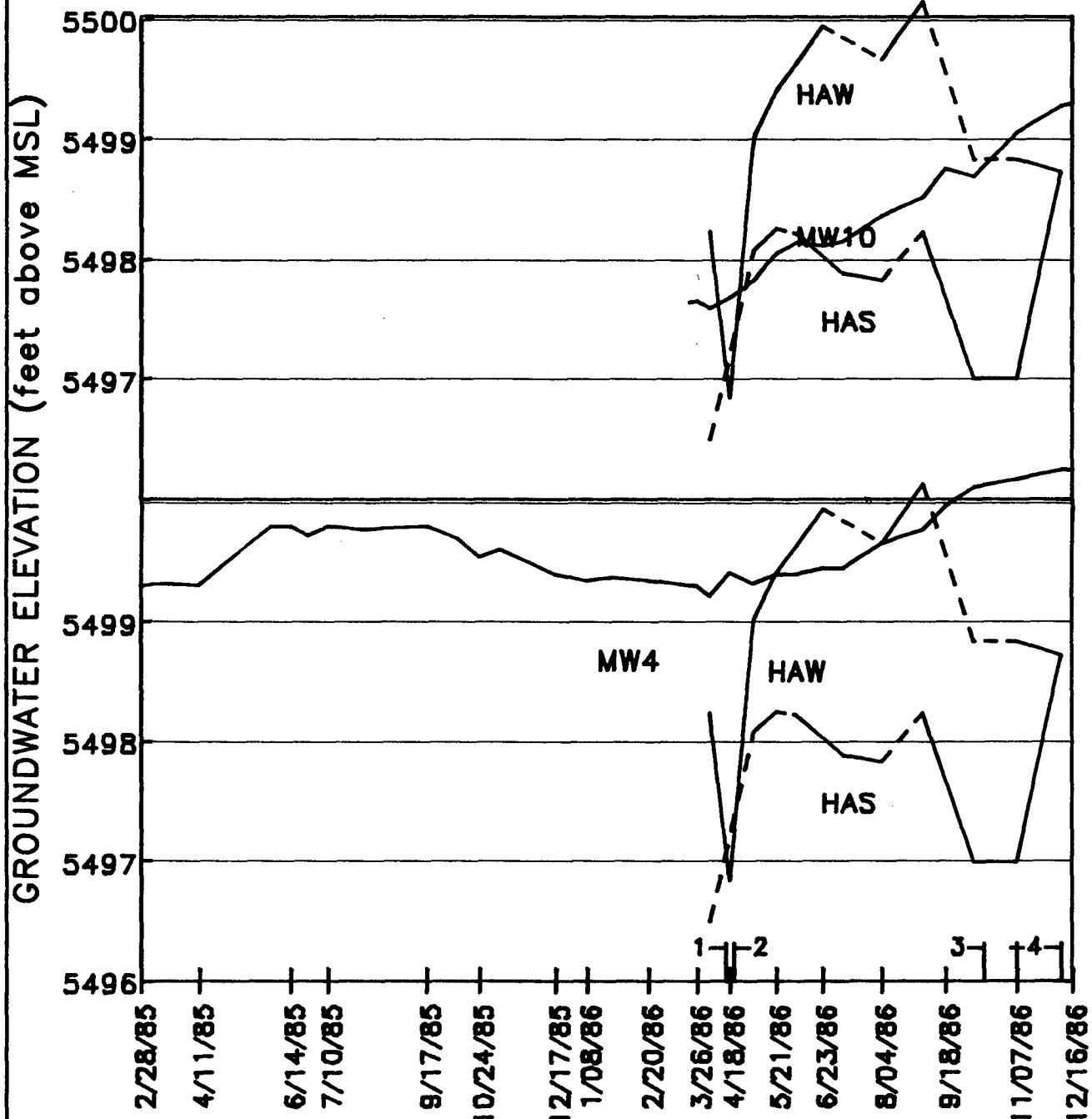
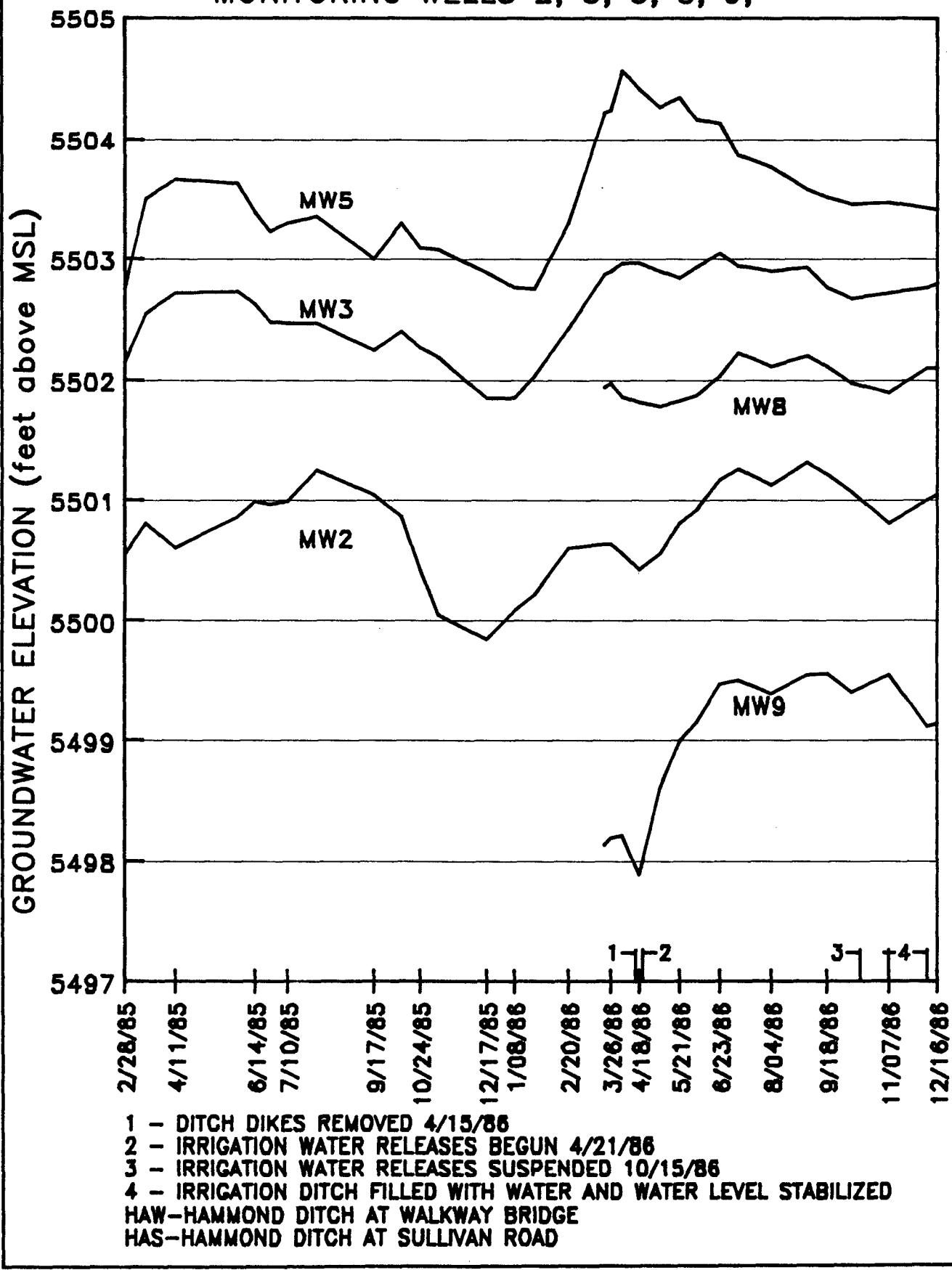


FIGURE 5.18
GROUNDWATER ELEVATIONS
BLOOMFIELD REFINING - BLOOMFIELD, NEW MEXICO
MONITORING WELLS 4, 10, HAW, HAS



still monitoring?

FIGURE 5.19
GROUNDWATER ELEVATIONS
BLOOMFIELD REFINING – BLOOMFIELD, NEW MEXICO
MONITORING WELLS 2, 3, 5, 8, 9,



wells, seasonal flows in Hammond Ditch affect these wells also. An additional year of water level data, which was also plotted, was available for wells 2, 3, and 5. The trend of the water levels in these wells is oscillatory, with periods of increasing water levels at the onset of irrigation season and corresponding decreases in water levels during nonirrigation periods.

Figure 5.20 was plotted to compare the relative water levels in wells 4 and 7 through time. Based on the performance of these wells, two trends become evident: Water levels in monitoring well 4 are consistently higher than those in monitoring well 7, and water level movements with time in monitoring well 7 appear to mimic closely those of well 4. This behavior suggests there is some small hydraulic influence at least to the superficial layers of the Nacimiento Formation. This influence is believed to be insignificant since water preferentially moves horizontally along the cobble zone/Nacimiento contact as evidenced by the seeps emerging at this contact along the bluffs of the San Juan River.

Another potentially significant facet of Figures 5.17 through 5.20 is the consistency at which reported water levels for December 1986 were approximately 1 foot higher than the water levels in January 1986. This consistency suggests that rainfall was more significant in 1986 than in 1985. Rainfall data were acquired by Bloomfield personnel. These data show that 1986 was a year of heavy rainfall. Table 5.2 shows the average total annual rainfall for the Bloomfield area is approximately 8.4 inches. This compares with 14.65 inches measured in 1986 - a 74 percent increase. Clarification of the unusual rainfall amounts in 1986 is shown in Table 5.3. This table shows that, for 9 months out of the 12, the 17-year monthly average rainfall was exceeded.

Ditch
may have
water all
winter
(from Hawley)

FIGURE 5.20
GROUNDWATER ELEVATIONS
BLOOMFIELD REFINING - BLOOMFIELD, NEW MEXICO
MONITORING WELLS 4 AND 7

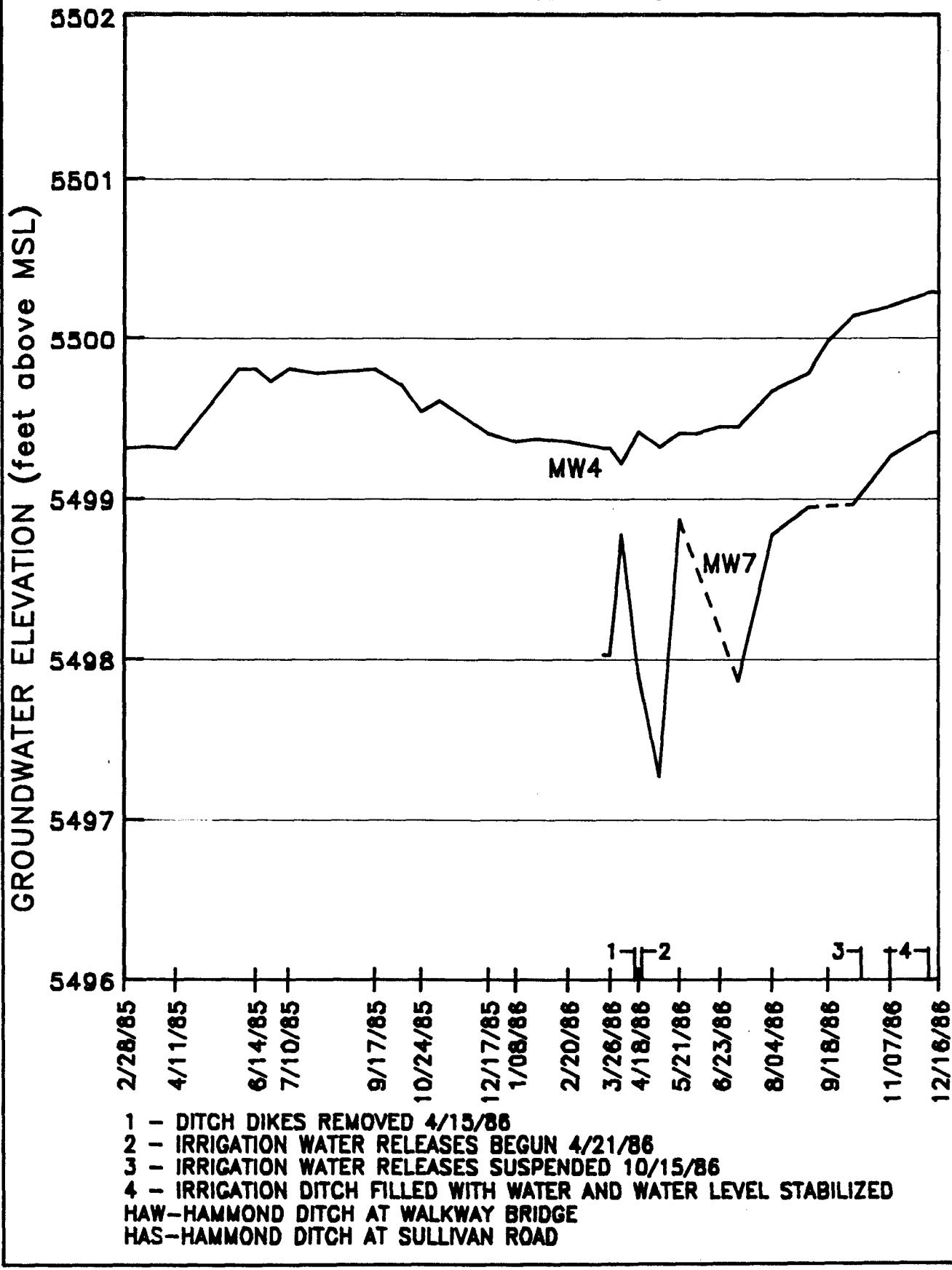


TABLE 5.2
HISTORICAL RAINFALL DATA FOR BLOOOMBFIELD REFINERY AREA
1969 THROUGH 1986*

| Month | 1969 | 1970 | 1971 | 1972 | 1973 | 1974 | 1975 | 1976 | 1977 | 1978 | 1979 | 1980 | 1981 | 1982 | 1983 | 1984 | 1985 | Avg. | 1986 |
|-----------|-------|------|------|------|------|------|------|------|------|-------|------|------|------|-------|------|------|------|------|-------|
| January | 0.85 | 0.06 | 0.10 | 0.03 | 0.78 | 1.1 | 0.51 | 0.06 | 0.42 | 0.9 | 0.88 | 1.45 | T | 0.32 | 0.94 | T | 0.39 | 0.52 | 0.11 |
| February | 0.31 | 0.07 | 0.09 | T | 0.12 | 0.13 | 0.61 | 0.16 | T | 0.84 | 0.79 | 0.70 | 0.30 | 0.77 | 0.69 | 0.12 | 0.13 | 0.45 | 0.77 |
| March | 0.21 | 0.47 | 0.03 | 0.07 | 1.62 | 0.01 | 1.52 | 0.00 | 0.00 | 1.21 | 0.46 | 0.63 | 1.76 | 1.18 | 1.84 | 0.54 | 1.74 | 1.28 | 0.51 |
| April | 0.3 | 0.38 | 0.71 | 0.08 | 1.34 | 0.2 | 0.18 | 0.1 | 0.07 | 0.14 | 0.2 | 0.25 | 0.21 | 0.67 | 0.31 | 1.00 | 1.76 | 0.70 | 0.97 |
| May | 1.44 | 0.11 | 0.42 | 0.02 | 0.63 | 0.02 | 0.75 | 0.41 | 0.79 | 0.96 | 0.58 | 0.25 | 1.05 | 0.82 | 0.13 | T | 0.29 | 0.42 | 0.13 |
| June | 1.00 | 0.81 | 0.08 | 0.18 | 0.93 | 0.09 | 0.11 | 0.09 | 0.04 | 0.69 | 0.43 | 0.07 | 0.16 | 0.00 | 0.35 | 0.67 | 0.01 | 0.21 | 0.81 |
| July | 0.69 | 0.68 | 0.34 | 0.04 | 0.22 | 1.09 | 0.68 | 0.42 | 1.07 | 0.01 | 1.4 | 0.08 | 1.34 | 1.27 | 1.67 | 0.62 | 1.38 | 1.06 | 4.10 |
| August | 0.41 | 0.02 | 1.72 | 1.54 | 0.81 | 0.12 | 0.24 | 0.8 | 1.41 | 0.18 | 0.4 | 0.89 | 0.35 | 2.78 | 0.72 | 1.64 | 0.43 | 1.14 | 0.93 |
| September | 2.01 | 2.68 | 1.00 | 0.57 | 1.49 | 0.37 | 0.8 | 1.41 | 0.78 | 1.35 | 0.08 | 1.05 | 0.69 | 1.50 | 0.53 | 0.45 | 1.31 | 0.92 | 2.18 |
| October | 2.00 | 0.48 | 1.15 | 2.57 | 0.25 | 2.12 | 0.14 | 0.09 | 0.3 | 1.4 | 1.71 | 0.84 | 0.89 | 0.16 | 0.52 | 1.13 | 1.21 | 0.79 | 0.65 |
| November | 0.78 | 0.46 | 0.21 | 0.79 | 0.3 | 0.48 | 0.22 | 0.01 | 0.62 | 2.24 | 0.43 | 0.02 | 0.36 | 0.92 | 0.91 | 0.23 | 0.52 | 0.49 | 2.73 |
| December | 0.29 | 0.2 | 0.16 | 0.97 | 0.31 | 0.38 | 0.2 | T | 0.63 | 0.59 | 0.23 | T | 0.03 | 0.76 | 0.67 | 0.87 | 0.22 | 0.43 | 0.76 |
| Total | 10.29 | 6.42 | 6.01 | 6.86 | 8.8 | 6.11 | 5.96 | 3.55 | 6.13 | 10.51 | 7.59 | 6.23 | 7.14 | 11.15 | 9.28 | 7.27 | 9.39 | 8.41 | 14.65 |

*Data for New Mexico State University Agricultural Center at Farmington, New Mexico.

TABLE 5.3

AVERAGE TEMPERATURE AND PRECIPITATION
FOR THE BLOOMFIELD REFINING AREA*

| Month | Average Precip. (inches) | Maximum Temperature (°F) | Minimum Temperature (°F) | Temperature Recorded | Extreme Maximum | | Year Recorded | Extreme Minimum |
|-----------|-----------------------------|-----------------------------|-----------------------------|----------------------|--------------------|------------------|-----------------------------|-----------------|
| | | | | | Year Recorded | Temperature (°F) | | |
| January | 0.47 | 39.6 | 18.4 | 61 | 1972 & 1975 | -18 | 1971 | |
| February | 0.30 | 47.0 | 23.1 | 68 | 1976 | -3 | 1982 | |
| March | 0.80 | 54.9 | 28.5 | 79 | 1978 | 6 | 1971 | |
| April | 0.52 | 64.6 | 34.4 | 83 | 1983 | 16 | 1979 | |
| May | 0.44 | 74.7 | 43.8 | 93 | 1974 & 1984 | 23 | 1975 | |
| June | 0.29 | 85.9 | 52.9 | 100 | 1981 | 36 | 1979, 1980, 1981, & 1983 | |
| July | 0.81 | 90.7 | 60.3 | 101 | 1971 | 43 | 1969 | |
| August | 0.84 | 88.0 | 59.3 | 99 | 1969, 1970, & 1983 | 41 | 1980 | |
| September | 1.04 | 79.9 | 51.2 | 95 | 1969 | 28 | 1971 | |
| October | 1.11 | 66.3 | 39.3 | 84 | 1975 & 1980 | 18 | 1971 | |
| November | 0.56 | 52.4 | 28.6 | 74 | 1977 | 1 | 1976 | |
| December | 0.41 | 42.6 | 20.5 | 65 | 1973 | -9 | 1978 | |
| Total | 7.59 | 65.6 | 38.4 | | | | | |

*Seventeen-year average weather conditions, New Mexico State University Agricultural Science Center at Farmington (1969-1985).

CHAPTER 6

ESTIMATION OF COBBLE LAYER AQUIFER CHARACTERISTICS

BACKGROUND

The EPA-approved work plan included specification of work to estimate the hydraulic characteristics of the upper cobble layer. Based on installation data of monitoring wells 1 through 6, wells 1, 2, and 4 were specified to be evaluated using slug tests. These wells were identified as most likely to yield appropriate information for estimation of aquifer characteristics and subsequently contaminant transport in the subsurface.

METHODOLOGY

On April 30, 1986, ES performed slug tests on wells 1, 2, and 4. These tests were performed by adding 16, 17, and 18 gallons to each well and measuring the time required for the casing water level to fall to its previously measured static level. The time required for the water level to return to static levels ranged from 2.75 minutes for well 2 to 7 minutes for well 4. Plots of the measured water level versus time were prepared and are presented as Figures 6.1 through 6.3.

Hydraulic conductivity was calculated for each of the wells. This calculation required investigation of the status of the screen in each well during the slug test measurements. Using well log information together with measurements of the depth of the well and the depth to water, factors affecting the calculation — the degree of siltation, penetration of impermeable Nacimiento Formation, and full or partial penetration of the aquifer — were taken into consideration. The saturated zone was also estimated from these measurements to enable calculation of a transmissivity. Averages of these numbers were calculated to derive "aquifer representative" values. Storage coefficient could not be obtained from slug test data but was estimated to be approximately the same as

FIGURE 6.1
SLUG TEST ANALYSIS
BLOOMFIELD REFINING
MW-1

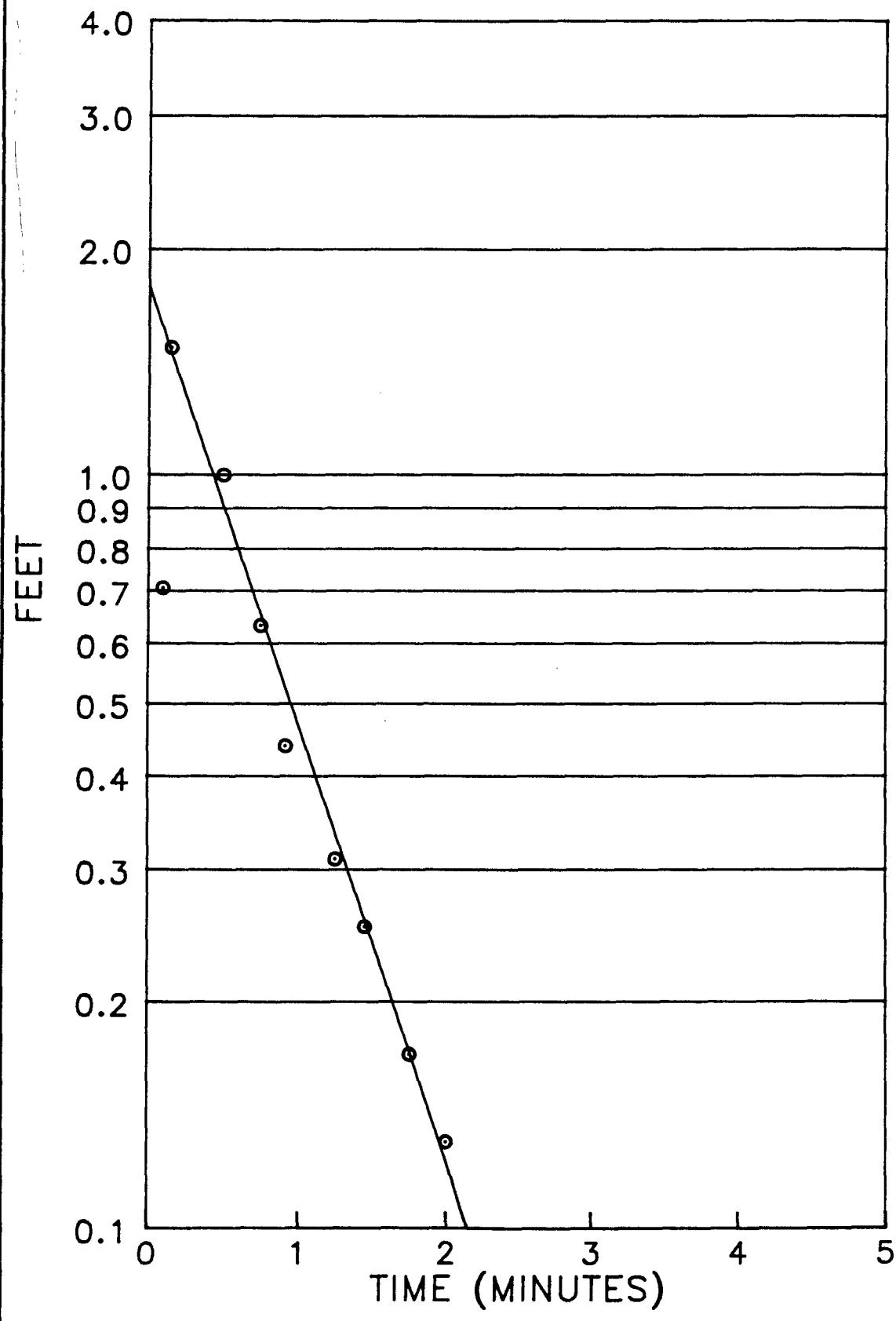


FIGURE 6.2
SLUG TEST ANALYSIS
BLOOMFIELD REFINING
MW-2

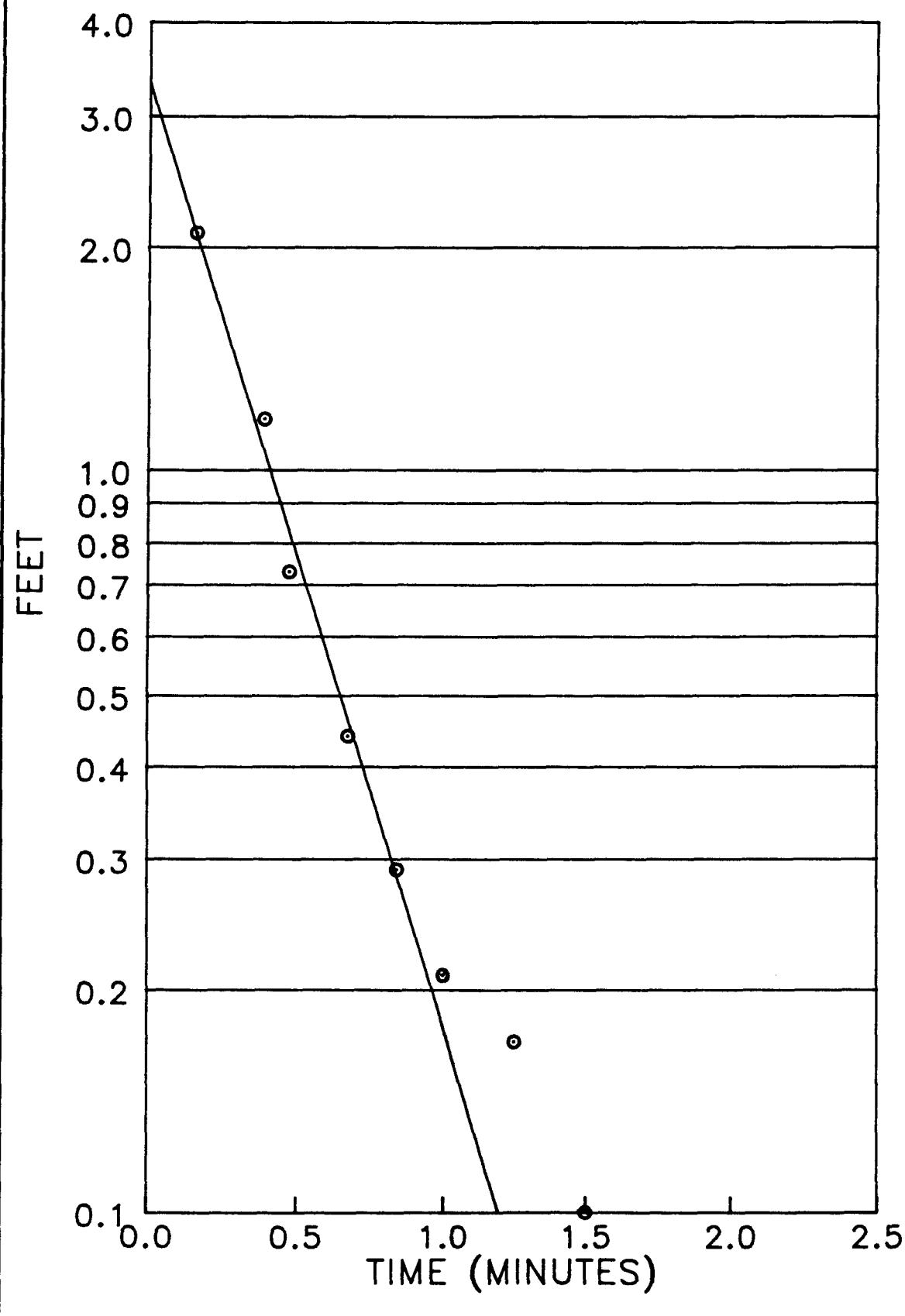
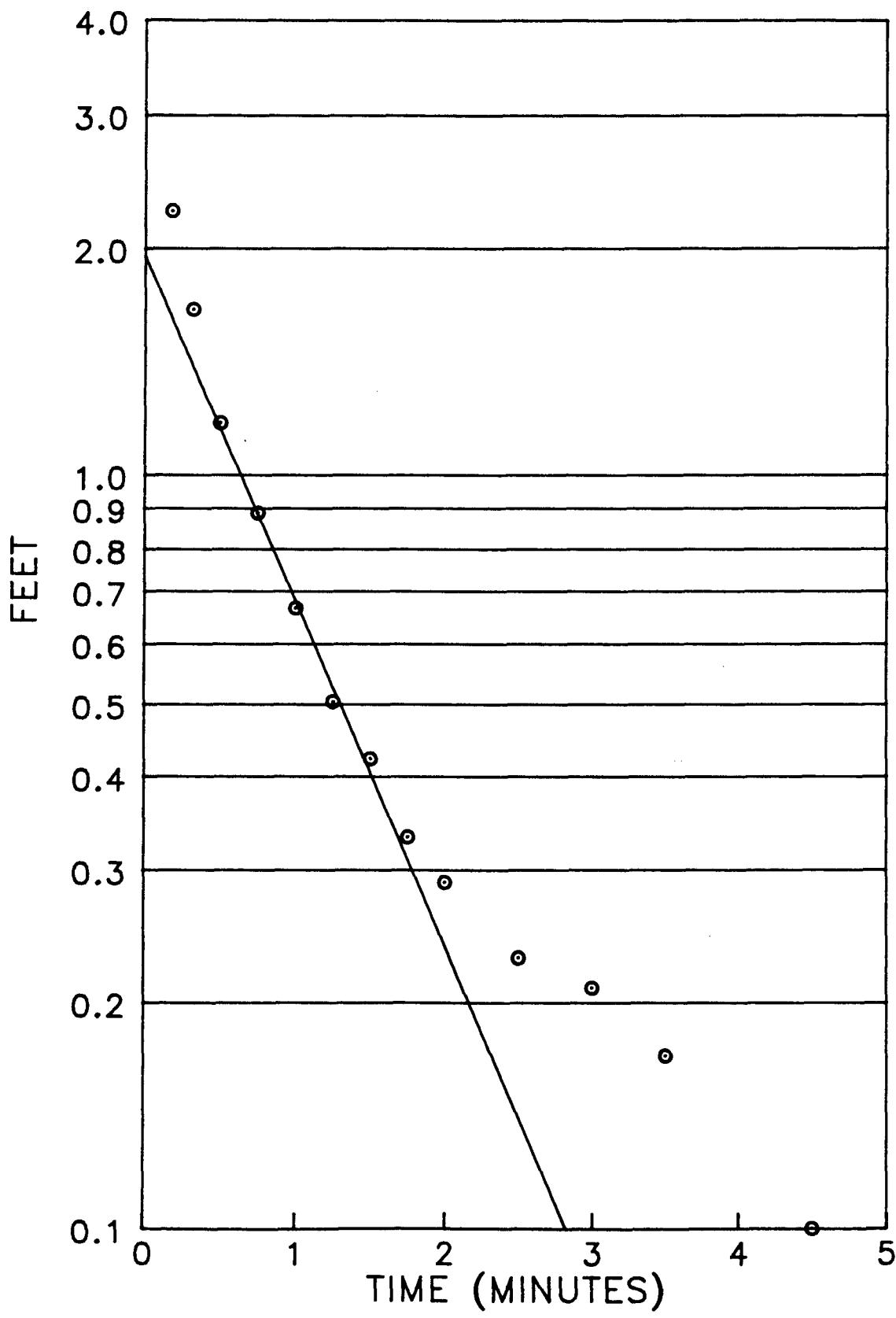


FIGURE 6.3
SLUG TEST ANALYSIS
BLOOMFIELD REFINING
MW-4



porosity. Porosity was assumed to be approximately 25 percent. A representative area of 200 feet square was assumed to provide a storage coefficient estimated to be 7.5×10^{-4} gallons per foot. ! Huh?

DISCUSSION

Based on the calculations in Table 6.1, the hydraulic conductivity of the cobble zone would appear to be an average value of 2.08×10^{-4} feet per second (6.34×10^{-3} centimeters per second). However, due to the shallow saturated thickness available, the transmissivity values are quite low and average 171 square feet per day. These factors, in combination with an average gradient estimated from groundwater contour maps in chapter 5 of 0.0025 and a width of approximately 2,500 feet (projection of the refinery perpendicular to flow), suggest a daily flux of approximately 8,500 gallons per day (gpd) or about 6 gpm.

TABLE 6.1
AQUIFER CHARACTERISTICS
OF THE COBBLE ZONE

| Monitoring Well Number | Estimated Saturated Thickness (feet) | Calculated Conductivity (ft/sec) | Transmissivity (square feet/day) |
|------------------------|--------------------------------------|--|----------------------------------|
| 1 | 10.6 | 1.65×10^{-4} | 151 |
| 2 | 9.2 | 3.30×10^{-4} | 262 |
| 4 | 9.0 | 1.29×10^{-4} | 100 |
| Average | 9.6 | 2.08×10^{-4} <i>185 ft/day</i> | 171 |

Ranges (from F. & Cherry): 60-70% Silty sand
40% Clean sand

$$Q = 2.08 \times 10^{-4} \text{ ft}^2/\text{sec} \times 0.0025 \times 2500 \text{ ft} \times 9.6 \text{ ft}$$

$$= 0.01258 \text{ ft}^3/\text{sec} = 8065 \text{ gpd} = 5.6 \text{ gpm}$$

CHAPTER 7

SURFACE WATER SAMPLING

BACKGROUND

Surface water sampling of Hammond Ditch and the San Juan River was required by the approved work plan. This sampling was specified to be accomplished during worst-case, low-flow conditions.

*Surf
Hammond
at 1*

Low-flow conditions were determined in the work plan to be at the beginning of the irrigation season for Hammond Ditch (generally mid-April). During this period, the potential for flushing hydrocarbons downstream is the greatest. BRC installs berms in Hammond Ditch at the end of the irrigation season to retain and remove hydrocarbons that may seep into the ditch. Just before resumption of irrigation, the berms are removed. The potential for downstream contamination is therefore greatest when upstream irrigation flows are released and flush the ditch.

Low-flow conditions in the San Juan River generally occur during winter months. Precipitation is generally low during winter. Throughout 1986, flows in the river were unusually high, and low-flow conditions did not occur. Maintenance activities for Navajo Dam (controlling San Juan River flows) in April and May 1987 required lowering the upstream reservoir approximately 31 feet from normal operating level. The lowering of the reservoir has resulted in unseasonably high discharge rates from the dam, which precluded low-flow sampling efforts. Therefore, an extension for sampling of the San Juan River has been requested to permit future sampling during representative low-flow conditions.

SAMPLING PROTOCOL

The Hammond Ditch was sampled by BRC staff on April 22, 1986 within 24 hours after initiation of irrigation flows in the ditch. Flows were observed to be intermittent during this period, so a second sampling effort was conducted on April 28, 1986 to assure acquisition of a representative data set.

All samples were collected consistent with the protocol stated in the approved work plan. As specified, depth-integrated samples were collected from a location just downstream of the refinery property south of Sullivan Road, and just downstream of the API wastewater ponds.

FINDINGS

Results from both Hammond Ditch sampling efforts are contained in Appendix B to this report. Tables 7.1 and 7.2 are summaries of compounds detected during the April 22, 1986 and April 28, 1986 sampling efforts, respectively. Table 7.1 shows that very low levels of pollutants were found during worst-case conditions. Table 7.2 shows that no priority pollutants were detected during the second survey, conducted after irrigation flows had been reestablished.

The results of the April 28, 1986 survey are consistent with the finding in Chapter 5 showing that Hammond Ditch irrigation flows create a hydraulic gradient directed towards the BRC site that minimizes the potential for groundwater transport into the ditch.

TABLE 7.1
SUMMARY OF DETECTED ORGANIC COMPOUNDS
FROM HAMMOND DITCH ON 4/22/86

| Laboratory Parameter | Hammond Ditch Near Sullivan Road Results (Downstream) | Hammond Ditch Near API Ponds Results (Upstream) | Detection Limits (mg/l) |
|----------------------|--|--|-------------------------------|
| | (mg/l) | (mg/l) | |
| Phenols | 0.002 | 0.002 | 0.001 |
| Benzene | 0.006 | | 0.001 |
| Toluene | 0.003 | | 0.001 |
| Anthracene | 0.006 | | 0.001 |
| Benzo(a)anthracene | 0.003 | | 0.001 |
| Chrysene | 0.005 | | 0.001 |
| Fluoranthene | | 0.001 | 0.001 |
| Naphthalene | 0.13 | | 0.001 |
| Phenanthrene | 0.008 | | 0.001 |
| Pyrene | 0.008 | | 0.001 |

TABLE 7.2
SUMMARY OF DETECTED ORGANIC COMPOUNDS
FROM HAMMOND DITCH ON 4/28/86

| Laboratory Parameter | Hammond Ditch Near Sullivan Road Results (Downstream) | Hammond Ditch Near API Ponds Results (Upstream) | Detection Limits (mg/l) |
|----------------------|--|--|-------------------------------|
| NONE DETECTED | | | |

CHAPTER 8

CONCLUSIONS

Each of the work elements specified in the approved work plan was designed to provide information that would clarify the investigation of subsurface hydrocarbons at the refinery. Based on execution of each of these work elements, the following conclusions are offered:

1. Subsurface petroleum hydrocarbons^(PRODUCT) appear to be located only within geologic units composed of sand, silt, clay, and cobbles. Subsurface hydrocarbon contamination may have been the result of old spills or leaks located within the vicinity of monitoring wells 4, 9, and 10.
2. Monitoring wells 1, 2, 3, 5, ⁽⁷⁾, and 8 show low levels of indicator pollutants.
3. Monitoring well 7 shows ^{see attached} insignificant levels of indicator pollutants. Well elevation data of wells 4 and 7 relative to changing water levels in Hammond Ditch suggest an ^{see attached} insignificant hydraulic influence between the screened portion of well 7 in the upper Nacimiento Formation and the cobble zone monitored by well 4.
4. No monitoring well has been established in a background area to clarify the significance or insignificance of metals and salts levels reported in water analyses. ^{see attached}
5. The ER profile maps, which provide an areal interpretation with respect to depth of discrete ER measurements, show existence of low ER contours located west of the solar evaporation ponds, underneath the process area and refined product storage areas, and offsite areas south and west of the intersection of Hammond Ditch with Sullivan Road. The presence of subsurface hydrocarbons has been indicated at monitoring wells 4, 9, and 10, but the areal extent and quantity of sources contributing to these wells are unknown. No information is available to explain the low ER contours predicted off site.

6. Hammond Ditch controls groundwater transport from the site towards the ditch during irrigation season. Elevated water levels in the ditch force groundwater that would otherwise enter the ditch to flow in a trough created by the intersecting gradients from the ditch and the refinery site.
7. Absence of water in Hammond Ditch removes the hydraulic impediment to flow towards Hammond Ditch. The extent of flow to Hammond Ditch and under or through Hammond Ditch is unknown.
8. Hammond Ditch surface water sampling showed that very low levels of pollutants are transmitted to Hammond Ditch during low-flow conditions. Nondetection of pollutants in the 4/28/86 sampling effort supports the conclusion that elevated Hammond Ditch water levels minimizes potential for transport towards the ditch.

CHAPTER 9

RECOMMENDATIONS

Based on the conclusions of the report, the following recommendations are offered:

1. During periods of nonirrigation flow, maintain elevated water levels in Hammond Ditch.
2. Pursue investigation of offsite conditions under auspices of the NMOCD.
3. Investigate information sources from which background water quality data exist or could be developed to reference the groundwater conditions beneath the BRC site.
4. Discontinue quality monitoring of all wells except 4, 10, and 9. Maintain quality monitoring of these wells on a quarterly basis for TOC.

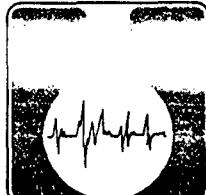
REFERENCES

- "A Work Plan for Monitoring, Testing, Analysis, and Reporting at Bloomfield Refinery." Engineering-Science, Inc. July 1985.
- "Discharge Plan for a Refinery Operated by Plateau, Inc. near Bloomfield, New Mexico." American Groundwater Consultants. March 24, 1984.
- "Guidance on Remedial Investigations Under CERCLA." U.S. EPA, EPA/540/6-85/002. June 1985.

APPENDIX A

MONITORING WELL SAMPLING RESULTS

FIRST QUARTER



ASSAIGAI ANALYTICAL LABORATORIES

SAMPLE DATE = 3/26/86

TO: Bloomfield Refinery
Attn: Chris Hawley
P.O. Box 159
Bloomfield, NM 87413

DATE: 14 May 1986
0502
Page 1 of 8

ANALYTE

SAMPLE ID/ ANALYTICAL RESULTS

| | MW 1 | MW 2 | MW 3 |
|-----------------|-------------|-------------|-------------|
| CN | <0.01 mg/l | <0.01 mg/l | <0.10 mg/l |
| Phenols | 0.009 mg/l | 0.063 mg/l | 0.006 mg/l |
| TOC | 18 mg/l | 18 mg/l | 29 mg/l |
| TDS | 2936 mg/l | 2796 mg/l | 4836 mg/l |
| Cl | 750 mg/l | 200 mg/l | 1500 mg/l |
| SO ₄ | 7.5 mg/l | 11.0 mg/l | 29.5 mg/l |
| Benzene | ND | ND | ND |
| Toluene | ND | ND | ND |
| Xylenes | ND | ND | ND |
| Ethylbenzene | ND | ND | ND |
| Sb | <0.01 mg/l | <0.01 mg/l | <0.01 mg/l |
| As | <0.050 mg/l | <0.050 mg/l | <0.050 mg/l |
| Be | <0.01 mg/l | <0.01 mg/l | <0.01 mg/l |
| Cd | 0.050 mg/l | 0.060 mg/l | 0.12 mg/l |
| Cr | <0.050 mg/l | <0.050 mg/l | <0.050 mg/l |
| Cu | <0.03 mg/l | <0.03 mg/l | <0.03 mg/l |
| Pb | 0.085 mg/l | 0.12 mg/l | 0.14 mg/l |
| Hg | <0.002 mg/l | 0.003 mg/l | 0.004 mg/l |
| Ni | 0.08 mg/l | 0.07 mg/l | 0.08 mg/l |
| Se | <0.010 mg/l | <0.010 mg/l | <0.010 mg/l |
| Ag | <0.050 mg/l | <0.050 mg/l | <0.050 mg/l |
| Tl | <0.01 mg/l | <0.01 mg/l | <0.01 mg/l |
| Zn | <0.01 mg/l | <0.01 mg/l | <0.01 mg/l |

TO: Bloomfield Refinery

DATE: 0502 revised

Page 2 of 8

SAMPLE DATE: 3/26/86

ANALYTE

SAMPLE ID/ ANALYTICAL RESULTS

| | MW 4 | MW 5 | MW 7 |
|--------------------------|-------------|-------------|-------------|
| CN | <0.01 mg/l | <0.01 mg/l | <0.10 mg/l |
| Phenols | 0.633 mg/l | 0.006 mg/l | <0.001 mg/l |
| TOC | 110 mg/l | 14 mg/l | 11 mg/l |
| TDS | 1868 mg/l | 3840 mg/l | 6076 mg/l |
| Cl | 500 mg/l | 1100 mg/l | 30 mg/l |
| SO 4 | 0.3 mg/l | 14.0 mg/l | 5.5 mg/l |
| Benzene | 11.8 mg/l | ND | 0.015 mg/l |
| Toluene | 7.5 mg/l | ND | 0.053 mg/l |
| Xylenes | ND | ND | ND |
| Ethylbenzene | 0.107 mg/l | ND | 0.007 mg/l |
| Sb | <0.01 mg/l | <0.01 mg/l | <0.01 mg/l |
| As | <0.050 mg/l | <0.050 mg/l | <0.050 mg/l |
| Be | <0.01 mg/l | <0.01 mg/l | <0.01 mg/l |
| Cd | 0.060 mg/l | 0.10 mg/l | 0.050 mg/l |
| Cr | <0.050 mg/l | <0.050 mg/l | <0.050 mg/l |
| Cu | <0.03 mg/l | <0.03 mg/l | <0.03 mg/l |
| Pb | 0.074 mg/l | 0.16 mg/l | <0.050 mg/l |
| Hg | 0.002 mg/l | <0.002 mg/l | <0.002 mg/l |
| Ni | 0.08 mg/l | 0.10 mg/l | 0.08 mg/l |
| Se | <0.010 mg/l | <0.010 mg/l | <0.010 mg/l |
| Ag | <0.050 mg/l | <0.050 mg/l | <0.050 mg/l |
| Tl | <0.01 mg/l | <0.01 mg/l | <0.01 mg/l |
| Zn | 0.012 mg/l | 0.012 mg/l | 0.018 mg/l |
| Acrolein | ND | | ND |
| Acrylonitrile | ND | | ND |
| Bromoform | ND | | ND |
| Carbon Tetrachloride | ND | | ND |
| Chlorobenzene | ND | | ND |
| Chlorodibromomethane | ND | | ND |
| Chloroethane | ND | | ND |
| 2-Chloroethylvinyl Ether | ND | | ND |
| Chloroform | ND | | ND |
| Dichlorogromomethane | ND | | ND |
| 1,1-Dichloroethane | ND | | ND |
| 1,2-Dichloroethane | ND | | ND |
| 1,1-Dichloroethylene | ND | | ND |
| 1,2-Dichloropropane | ND | | ND |
| 1,3-Dichloropropylene | ND | | ND |

SAMPLE DATE: 3/20/00

TO: Bloomfield Refinery

0502

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ANALYTE

SAMPLE ID/ ANALYTICAL RESULTS

MW 4

MW 5

MW 7

| | | |
|-----------------------------|------------|------------|
| Methyl Bromide | ND | ND |
| Methyl Chloride | ND | ND |
| Methylene Chloride | ND | ND |
| 1,1,2,2-Tetrachloroethane | ND | ND |
| Tetrachloroethylene | ND | ND |
| 1,2-Transdichloroethylene | ND | ND |
| 1,1,1-Trichloroethane | ND | ND |
| 1,1,2-Trichloroethane | ND | ND |
| Trichloroethylene | ND | ND |
| Vinyl Chloride | ND | ND |
| Acid Compounds | | |
| 2-chlorophenol | ND | ND |
| 2,4-dichlorophenol | 0.200 mg/l | ND |
| 2,4-dimethylphenol | ND | ND |
| 4,6-dinitro-o-cresol | 0.100 mg/l | 0.013 mg/l |
| 2,4-dinitrophenol | 0.050 mg/l | ND |
| 2-nitrophenol | ND | ND |
| 4-nitrophenol | 0.090 mg/l | ND |
| p-chloro-m-cresol | ND | ND |
| pentachlorophenol | ND | ND |
| Phenol | 0.202 mg/l | ND |
| 2,4,6-trichlorophenol | ND | ND |
| Base Neutrals | | |
| Acenaphthene | 0.044 mg/l | ND |
| Acenaphthylene | ND | ND |
| Anthracene | ND | ND |
| Benzidine | ND | ND |
| Benzo(a)anthracene | ND | ND |
| Benzo(a)pyrene | ND | ND |
| 3,4-benzofluoranthene | ND | ND |
| Benzo(ghi)perylene | ND | ND |
| Benzo(k)fluoranthene | ND | ND |
| Bis(2-chloroethoxy)methane | ND | ND |
| Bis(2-chloroethyl)ether | ND | ND |
| Bis(2-chloroisopropyl)ether | ND | ND |
| Bis(2-ethylhexyl)phthalate | ND | ND |
| 4-bromophenyl phenyl ether | ND | ND |
| Butylbenzyl phthalate | ND | ND |
| 2-chloronaphthalene | ND | ND |
| 4-chlorophenyl phenyl ether | ND | ND |
| Chrysene | ND | ND |

SAMPLE DATE : 3/26/86

TO: Bloomfield Refinery

0502

Page 4 of 8

ANALYTE

SAMPLE ID/ ANALYTICAL RESULTS

MW 4

MW 5

MW 7

| | | |
|---------------------------|------------|----|
| Dibenzo(a,h)anthracene | ND | ND |
| 1,2-Dichlorobenzene | ND | ND |
| 1,3-Dichlorobenzene | ND | ND |
| 1,4-Dichlorobenzene | ND | ND |
| 3,3-Dichlorobenzidine | ND | ND |
| Diethyl phthalate | ND | ND |
| Dimethyl phthalate | ND | ND |
| Din-n-butyl phthalate | ND | ND |
| 2,4-dinitrotoluene | ND | ND |
| 2,6-dinitrotoluene | ND | ND |
| Di-n-octyl phthalate | ND | ND |
| 1,2-diphenylhydrazine | ND | ND |
| Fluoranthene | ND | ND |
| Fluorene | 0.150 mg/l | ND |
| Hexachlorobenzene | ND | ND |
| Hexachlorobutadiene | ND | ND |
| Hexachlorocyclopentadiene | ND | ND |
| Hexachloroethane | ND | ND |
| Indeno(1,2,3-cd)pyrene | ND | ND |
| Isophorone | ND | ND |
| Naphthalene | 0.036 mg/l | ND |
| Nitrobenzene | ND | ND |
| N-nitrosodimethylamine | ND | ND |
| N-nitrosodi-n-propylamine | ND | ND |
| N-nitrosodiphenylamine | ND | ND |
| Phenanthrene | ND | ND |
| Pyrene | 0.166 mg/l | ND |
| 1,2,4-trichlorobenzene | ND | ND |

SAMPLE DATE: 3/6/86

TO: Bloomfield Refinery

DATE: 0502

Page 5 of 8

ANALYTE

SAMPLE ID/ ANALYTICAL RESULTS

| | MW 8 | MW 9 | MW 10 |
|--------------------------|-------------|-------------|-------------|
| CN | <0.01 mg/l | <0.01 mg/l | <0.10 mg/l |
| Phenols | <0.001 mg/l | 0.304 mg/l | 0.147 mg/l |
| TOC ^X | 5 mg/l | 143 mg/l | 34 mg/l |
| TDS | 806 mg/l | 2360 mg/l | 1546 mg/l |
| Cl | 160 mg/l | 149 mg/l | 245 mg/l |
| SO ₄ | 4.0 mg/l | 13.0 mg/l | 5.3 mg/l |
| Benzene | ND | 7.4 mg/l | 0.093 mg/l |
| Toluene | ND | 6.3 mg/l | ND |
| Xylenes | ND | ND | ND |
| Ethylbenzene | 0.107 mg/l | 3.2 mg/l | ND |
| Sb ^X | <0.01 mg/l | <0.01 mg/l | <0.01 mg/l |
| As | <0.050 mg/l | <0.050 mg/l | <0.050 mg/l |
| Be ^X | <0.01 mg/l | <0.01 mg/l | <0.01 mg/l |
| Cd | 0.010 mg/l | 0.010 mg/l | 0.020 mg/l |
| Cr | <0.050 mg/l | <0.050 mg/l | <0.050 mg/l |
| Cu ^X | <0.03 mg/l | <0.03 mg/l | <0.03 mg/l |
| Pb | <0.050 mg/l | <0.050 mg/l | <0.050 mg/l |
| Hg ^X | <0.002 mg/l | <0.002 mg/l | <0.002 mg/l |
| Ni | <0.06 mg/l | 0.30 mg/l | 0.08 mg/l |
| Se | <0.010 mg/l | <0.010 mg/l | <0.010 mg/l |
| Ag ^X | <0.050 mg/l | <0.050 mg/l | <0.050 mg/l |
| Tl ^X | <0.01 mg/l | <0.01 mg/l | <0.01 mg/l |
| Zn | <0.01 mg/l | 0.012 mg/l | <0.01 mg/l |
| Acrolein | ND | ND | ND |
| Acrylonitrile | ND | ND | ND |
| Bromoform | ND | ND | ND |
| Carbon Tetrachloride | ND | ND | ND |
| Chlorobenzene | ND | ND | ND |
| Chlorodibromomethane | ND | ND | ND |
| Chloroethane | ND | ND | ND |
| 2-Chloroethylvinyl Ether | ND | ND | ND |
| Chloroform | ND | ND | ND |
| Dichlorogromomethane | ND | ND | ND |
| 1,1-Dichloroethane | ND | ND | ND |
| 1,2-Dichloroethane | ND | ND | ND |
| 1,1-Dichloroethylene | ND | ND | ND |
| 1,2-Dichloropropane | ND | ND | ND |
| 1,3-Dichloropropylene | ND | ND | ND |

SAMPLE DATE - 3/26/96

TO: Bloomfield Refinery

0502

Page 6 of 8

ANALYTE

SAMPLE ID/ ANALYTICAL RESULTS

MW 8 MW 9 MW 10

| | | | |
|-----------------------------|----|------------|------------|
| Methyl Bromide | ND | ND | ND |
| Methyl Chloride | ND | ND | ND |
| Methylene Chloride | ND | ND | ND |
| 1,1,2,2-Tetrachloroethane | ND | ND | ND |
| Tetrachloroethylene | ND | ND | ND |
| 1,2-Transdichloroethylene | ND | ND | ND |
| 1,1,1-Trichloroethane | ND | ND | ND |
| 1,1,2-Trichloroethane | ND | ND | ND |
| Trichloroethylene | ND | ND | ND |
| Vinyl Chloride | ND | ND | ND |
| Acid Compounds | | | |
| 2-chlorophenol | ND | ND | ND |
| 2,4-dichlorophenol | ND | ND | ND |
| 2,4-dimethylphenol | ND | 0.160 mg/l | 0.025 mg/l |
| 4,6-dinitro-o-cresol | ND | ND | 0.020 mg/l |
| 2,4-dinitrophenol | ND | ND | ND |
| 2-nitrophenol | ND | ND | ND |
| 4-nitrophenol | ND | ND | ND |
| p-chloro-m-cresol | ND | ND | ND |
| pentachlorophenol | ND | ND | ND |
| Phenol | ND | 0.149 mg/l | 0.090 mg/l |
| 2,4,6-trichlorophenol | ND | ND | ND |
| Base Neutrals | | | |
| Acenaphthene | ND | ND | ND |
| Acenaphthylene | ND | ND | ND |
| Anthracene | ND | ND | 0.039 mg/l |
| Benzidine | ND | ND | ND |
| Benzo(a)anthracene | ND | ND | ND |
| Benzo(a)pyrene | ND | ND | ND |
| 3,4-benzofluoranthene | ND | ND | ND |
| Benzo(ghi)perylene | ND | ND | ND |
| Benzo(k)fluoranthene | ND | ND | ND |
| Bis(2-chloroethoxy)methane | ND | ND | ND |
| Bis(2-chloroethyl)ether | ND | ND | ND |
| Bis(2-chloroisopropyl)ether | ND | ND | ND |
| Bis(2-ethylhexyl)phthalate | ND | ND | ND |
| 4-bromophenyl phenyl ether | ND | ND | ND |
| Butylbenzyl phthalate | ND | ND | ND |
| 2-chloronaphthalene | ND | ND | ND |
| 4-chlorophenyl phenyl ether | ND | ND | ND |
| Chrysene | ND | ND | ND |

SAMPLE DATE : 3/26/86

TO: Bloomfield Refinery

0502

Page 7 of 8

ANALYTE

SAMPLE ID/ ANALYTICAL RESULTS

| | MW 8 | MW 9 | MW 10 |
|----------------------------|------|------------|------------|
| Dibenzo(a,h)anthracene | ND | ND | ND |
| 1,2-Dichlorobenzene | ND | ND | ND |
| 1,3-Dichlorobenzene | ND | ND | ND |
| 1,4-Dichlorobenzene | ND | ND | ND |
| 3,3-Dichlorobenzidine | ND | ND | ND |
| Diethyl phthalate | ND | ND | ND |
| Dimethyl phthalate | ND | ND | ND |
| Din-n-butyl phthalate | ND | ND | ND |
| 2,4-dinitrotoluene | ND | ND | ND |
| 2,6-dinitrotoluene | ND | ND | ND |
| Di-n-octyl phthalate | ND | ND | ND |
| 1,2-diphenylhydrazine | ND | ND | ND |
| Fluoranthene | ND | ND | 0.034 mg/l |
| Fluorene | ND | 0.012 mg/l | 0.033 mg/l |
| Hexachlorobenzene | ND | ND | ND |
| Hexachlorobutadiene | ND | ND | ND |
| Hexachlorocyclopentadiene | ND | ND | ND |
| Hexachloroethane | ND | ND | ND |
| Indeno(1,2,3-cd)pyrene | ND | ND | ND |
| Isophorone | ND | ND | ND |
| Naphthalene | ND | ND | ND |
| Nitrobenzene | ND | ND | ND |
| N-nitrosodimethylamine | ND | ND | ND |
| N-nitrosodie-n-propylamine | ND | ND | ND |
| N-nitrosodiphenylamine | ND | ND | ND |
| Phenanthrene | ND | ND | ND |
| Pyrene | ND | ND | 0.030 mg/l |
| 1,2,4-trichlorobenzene | ND | ND | ND |

ND = None Detected

REFERENCE: "Test Methods for Evaluating Solid Waste, Physical/ Chemical Methods, USEPA, SW 846, EMSL-Cincinnati, 1982.

SAMPLE DATE: 3/26/86

TO: Bloomfield Refinery

0502 revised
Page 8 Of 8

NOMINAL DETECTION LIMITS

| | |
|--------------|------------|
| CN | 0.01 mg/l |
| Phenols | 0.001 mg/l |
| TOC | 0.1 mg/l |
| TDS | 1 mg/l |
| Cl | 1.0 mg/l |
| SO 4 | 1.0 mg/l |
| Benzene | 0.005 mg/l |
| Toluene | 0.005 mg/l |
| Xylenes | 0.005 mg/l |
| Ethylbenzene | 0.005 mg/l |
| Sb | 0.01 mg/l |
| As | 0.050 mg/l |
| Be | 0.01 mg/l |
| Cd | 0.002 mg/l |
| Cr | 0.050 mg/l |
| Cu | 0.03 mg/l |
| Pb | 0.001 mg/l |
| Hg | 0.002 mg/l |
| Ni | 0.01 mg/l |
| Se | 0.010 mg/l |
| Ag | 0.050 mg/l |
| Tl | 0.01 mg/l |
| Zn | 0.01 mg/l |

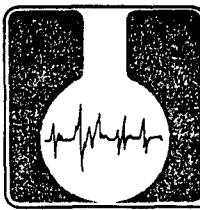
Detection limits for Volatiles, Acid Compounds, and Base/Neutrals are
all 0.001 mg/l

An invoice for services is enclosed. Thank you for contacting Assaigai
Laboratories.

Sincerely,

Jennifer V. Smith, Ph.D.
Laboratory Director

SECOND QUARTER



ASSAIGAI ANALYTICAL LABORATORIES

TO: Bloomfield Refinery
Attn: Chris Hawley
PO Box 159
Bloomfield, NM 87413

DATE: 23 July 1986
1030

SAMPLE ID: MW - 1

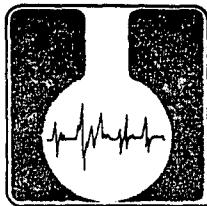
ANALYTE

ANALYTICAL RESULTS

| | |
|--------------|-----------------------|
| CN | 0.1 mg/l |
| TDS | 2960 mg/l |
| Cl | 994.7 mg/l |
| SO 4 | 630 mg/l |
| Phenols | 0.017 mg/l |
| TOC | 24 mg/l |
| Sb | <0.01 mg/l |
| As | 0.077 mg/l |
| Be | <0.01 mg/l |
| Cd | <0.010 mg/l |
| Cr | <0.050 mg/l |
| Cu | <0.050 mg/l duplicate |
| Pb | <0.03 mg/l |
| Hg | 0.065 mg/l |
| Ni | <0.002 mg/l |
| Se | <0.06 mg/l |
| Ag | 0.035 mg/l |
| Tl | <0.050 mg/l |
| Zn | <0.01 mg/l |
| Benzene | 0.020 mg/l |
| Toluene | ND |
| Xylenes | ND |
| Ethylbenzene | ND |
| Ba | <0.01 mg/l |
| Fe | <0.04 mg/l |
| Mn | 0.25 mg/l |
| Al | 2.07 mg/l |
| B | <0.01 mg/l |
| Co | <0.05 mg/l |
| Mo | <0.01 mg/l |
| F | 0.54 mg/l |
| No 3 as N | 0.1 mg/l |
| 1,2-DCE | ND |
| 1,1-DCE | ND |
| 1,1,2,2-TCE | ND |
| 1,1,2-TCE | ND |

Field by C4 6/23/86

pH 7.25
Conductivity 4600



ASSAIGAI ANALYTICAL LABORATORIES

TO: Bloomfield Refinery
Attn: Chris Hawley
PO Box 159
Bloomfield, NM 87413

DATE: 23 July 1986
1030

SAMPLE ID: MW - 2

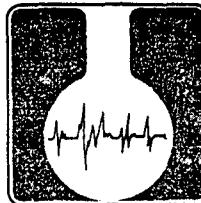
ANALYTE

ANALYTICAL RESULTS

| | |
|--------------|-------------|
| CN | 0.1 mg/l |
| TDS | 3650 mg/l |
| Cl | 1204.6 mg/l |
| SO 4 | 1750 mg/l |
| Phenols | 0.023 mg/l |
| TOC | 27 mg/l |
| Sb | <0.01 mg/l |
| As | 0.094 mg/l |
| Be | <0.01 mg/l |
| Cd | <0.010 mg/l |
| Cr | <0.050 mg/l |
| Cu | <0.03 mg/l |
| Pb | <0.05 mg/l |
| Hg | <0.002 mg/l |
| Ni | <0.06 mg/l |
| Se | 0.070 mg/l |
| Ag | <0.050 mg/l |
| Tl | <0.01 mg/l |
| Zn | 0.020 mg/l |
| Benzene | ND |
| Toluene | ND |
| Xylenes | ND |
| Ethylbenzene | ND |

Field by C4 6/23/86

pH 7.17
Conductivity 5400



ASSAIGAI ANALYTICAL LABORATORIES

TO: Bloomfield Refinery
Attn: Chris Hawley
PO Box 159
Bloomfield, NM 87413

DATE: 23 July 1986
1030

SAMPLE ID: MW - 3

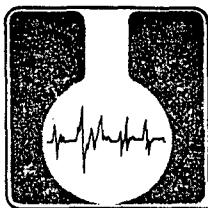
ANALYTE

ANALYTICAL RESULTS

| | |
|--------------|-------------|
| CN | 0.25 mg/l |
| TDS | 5362 mg/l |
| Cl | 1584 mg/l |
| SO 4 | 1950 mg/l |
| Phenols | 0.006 mg/l |
| TOC | 17 mg/l |
| Sb | <0.01 mg/l |
| As | 0.15 mg/l |
| Be | <0.01 mg/l |
| Cd | 0.015 mg/l |
| Cr | <0.050 mg/l |
| Cu | <0.03 mg/l |
| Pb | 0.070 mg/l |
| Hg | <0.002 mg/l |
| Ni | 0.08 mg/l |
| Se | 0.10 mg/l |
| Ag | <0.050 mg/l |
| Tl | <0.01 mg/l |
| Zn | 0.018 mg/l |
| Benzene | ND |
| Toluene | 0.003 mg/l |
| Xylenes | 0.030 mg/l |
| Ethylbenzene | ND |

Field by Cht 6/23/86

pH 7.10
Conductivity 6900



ASSAIGAI ANALYTICAL LABORATORIES

TO: Bloomfield Refinery
Attn: Chris Hawley
PO Box 159
Bloomfield, NM 87413

DATE: 23 July 1986
1030

SAMPLE ID: MW - 4

| ANALYTE | ANALYTICAL RESULTS |
|--------------------------|--------------------|
| CN | 0.5 mg/l |
| TDS | 2266 mg/l |
| Cl | 989.7 mg/l |
| SO 4 | 12.5 mg/l |
| Phenols | 0.430 mg/l |
| TOC | 130 mg/l |
| Sb | <0.10 mg/l |
| As | 0.070 mg/l |
| Be | <0.1 mg/l |
| Cd | <0.010 mg/l |
| Cr | <0.050 mg/l |
| Cu | <0.03 mg/l |
| Pb | 0.066 mg/l |
| Hg | <0.002 mg/l |
| Ni | <0.06 mg/l |
| Se | 0.080 mg/l |
| Ag | <0.050 mg/l |
| Tl | <0.1 mg/l |
| Zn | 0.019 mg/l |
| Volatiles | |
| Acrolein | ND |
| Acrylonitrile | ND |
| Benzene | 3.1 mg/l |
| Bromoform | ND |
| Carbon Tetrachloride | ND |
| Chlorobenzene | ND |
| Chlorodibromomethane | ND |
| Chloroethane | ND |
| 2-Chloroethylvinyl ether | ND |
| Chloroform | ND |
| Dichlorobromomethane | ND |
| 1,1-Dichloroethane | ND |
| 1,2-Dichloroethane | ND |
| 1,1-Dichloroethylene | ND |
| 1,2-Dichloropropane | ND |
| 1,2-Dichloropropylene | ND |
| Ethylbenzene | 0.070 mg/l |

Field by CL 6/4/86

pH 6.85

Conductivity 3800

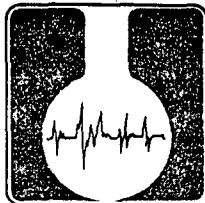
SAMPLE ID: MW - 4

| ANALYTE | ANALYTICAL RESULTS |
|-----------------------------|--------------------|
| Methyl Bromide | ND |
| Methyl Chloride | ND |
| Methylene Chloride | ND |
| 1,1,2,2-Tetrachloroethane | ND |
| Tetrachloroethylene | ND |
| Toluene | 0.290 mg/l |
| 1,2-Transdichloroethylene | ND |
| 1,1,1-Trichloroethane | ND |
| 1,1,2-Trichloroethane | ND |
| Trichloroethylene | ND |
| Vinyl Chloride | ND |
| Acid Compounds | |
| 2-Chlorophenol | ND |
| 2,4-Dichlorophenol | ND |
| 2,4-Dimethylphenol | 0.058 mg/l |
| 4,6-Dinitro-o-cresol | ND |
| 2,4-Dinitrophenol | ND |
| 2-Nitrophenol | 0.108 mg/l |
| 4-Nitrophenol | 0.302 mg/l |
| P-chloro-m-cresol | ND |
| pentachlorophenol | ND |
| Phenol | ND |
| 2,4,6-Trichlorophenol | ND |
| Base Neutrals | |
| Acenaphthene | ND |
| Acenaphthylene | ND |
| Anthracene | ND |
| Benzidine | ND |
| Benzo(a)anthracene | 0.016 mg/l |
| Benzo(a)pyrene | ND |
| 3,4-Benzofluoranthene | ND |
| Benzo(g,h,i)perylene | ND |
| Benzo(k)fluoranthene | ND |
| Bis(2-chloroethoxy)methane | ND |
| Bis(2-chloroethyl)ether | ND |
| Bis(2-chloroisopropyl)ether | ND |
| Bis(2-ethylhexyl)phthalate | ND |
| 4-Bromophenyl phenyl ether | ND |
| Butylbenzyl phthalate | ND |
| 2-Chloronaphthalene | ND |
| 4-Chlorophenyl phenyl ether | ND |
| Chrysene | 0.023 mg/l |

SAMPLE ID: MW - 4

| ANALYTE | ANALYTICAL RESULTS |
|----------------------------|--------------------|
| Dibenzo(a,h)anthracene | ND |
| 1,2-Dichlorobenzene | ND |
| 1,3-Dichlorobenzene | ND |
| 1,4-Dichlorobenzene | ND |
| 3,3-Dichlorobenzidine | ND |
| Diethyl phthalate | ND |
| Dimethyl phthalate | ND |
| Din-n-butyl phthalate | ND |
| 2,4-Dinitrotoluene | ND |
| 2,6-Dinitrotoluene | ND |
| Di-n-octyl phthalate | ND |
| 1,2-Diphenylhydrazine | ND |
| Fluoranthene | ND |
| Fluorene | ND |
| Hexachlorobenzene | ND |
| Hexachlorobutadiene | ND |
| Hexachlorocyclopentadiene | ND |
| Hexachloroethane | ND |
| Indeno(1,2,3-cd)pyrene | ND |
| Isophorone | ND |
| Naphthalene | 0.019 mg/l |
| Nitrobenzene | ND |
| N-nitrosodimethylamine | ND |
| N-nitrosodie-n-propylamine | ND |
| N-nitrosodiphenylamine | ND |
| Phenanthrene | ND |
| Pyrene | ND |
| 1,2,4-Trichlorobenzene | ND |
| Ba | 3.54 mg/l |
| Fe | 12.0 mg/l |
| Mn | 3.5 mg/l |
| Al | 1.93 mg/l |
| B | <0.01 mg/l |
| Co | <0.05 mg/l |
| Mo | <0.01 mg/l |
| F | 0.21 mg/l |
| NO ₃ as N | <0.01 mg/l |

ND = None Detected



ASSAIGAI ANALYTICAL LABORATORIES

TO: Bloomfield Refinery
Attn: Chris Hawley
PO Box 159
Bloomfield, NM 87413

DATE: 23 July 1986
1030

SAMPLE ID: MW - 5

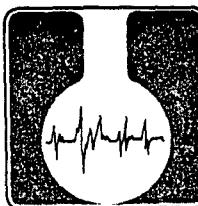
ANALYTE

ANALYTICAL RESULTS

| | |
|--------------|-------------|
| CN | 0.2 mg/l |
| TDS | 3778 mg/l |
| Cl | 1339.6 mg/l |
| SO 4 | 1800 mg/l |
| Phenols | 0.007 mg/l |
| TOC | 21 mg/l |
| Sb | <0.01 mg/l |
| As | 0.087 mg/l |
| Be | <0.01 mg/l |
| Cd | <0.010 mg/l |
| Cr | <0.050 mg/l |
| Cu | <0.03 mg/l |
| Pb | 0.055 mg/l |
| Hg | <0.002 mg/l |
| Ni | <0.06 mg/l |
| Se | 0.071 mg/l |
| Ag | <0.050 mg/l |
| Tl | <0.01 mg/l |
| Zn | 0.02 mg/l |
| Benzene | ND |
| Toluene | ND |
| Xylenes | ND |
| Ethylbenzene | ND |
| Ba | <0.01 mg/l |
| Fe | 0.05 mg/l |
| Mn | 0.025 mg/l |
| Al | 2.75 mg/l |
| B | <0.01 mg/l |
| Co | <0.05 mg/l |
| Mo | <0.01 mg/l |
| F | 0.30 mg/l |
| No 3 as N | 12.5 mg/l |
| 1,2-DCE | ND |
| 1,1-DCE | ND |
| 1,1,2,2-TCE | ND |
| 1,1,2-TCE | ND |

Field by C14 6/23/86

pH 7.18
Conductivity 5400



ASSAIGAI ANALYTICAL LABORATORIES

TO: Bloomfield Refinery
Attn: Chris Hawley
PO Box 159
Bloomfield, NM 87413

DATE: 23 July 1986
1030

SAMPLE ID: MW - 7

ANALYTE

ANALYTICAL RESULTS

| | |
|--------------------------|-------------|
| CN | 0.25 mg/l |
| TDS | 6406 mg/l |
| Cl | 79.9 mg/l |
| SO 4 | 2400 mg/l |
| Phenols | 0.006 mg/l |
| TOC | 4 mg/l |
| Sb | <0.01 mg/l |
| As | 0.36 mg/l |
| Be | <0.01 mg/l |
| Cd | 0.030 mg/l |
| Cr | 0.052 mg/l |
| Cu | <0.03 mg/l |
| Pb | 0.24 mg/l |
| Hg | <0.002 mg/l |
| Ni | 0.07 mg/l |
| Se | 0.65 mg/l |
| Ag | 0.060 mg/l |
| Tl | <0.01 mg/l |
| Zn | 0.016 mg/l |
| Volatiles | |
| Acrolein | ND |
| Acrylonitrile | ND |
| Benzene | ND |
| Bromoform | ND |
| Carbon Tetrachloride | ND |
| Chlorobenzene | ND |
| Chlorodibromomethane | ND |
| Chloroethane | ND |
| 2-Chloroethylvinyl ether | ND |
| Chloroform | ND |
| Dichlorobromomethane | ND |
| 1,1-Dichloroethane | ND |
| 1,2-Dichloroethane | ND |
| 1,1-Dichloroethylene | ND |
| 1,2-Dichloropropane | ND |
| 1,2-Dichloropropylene | ND |
| Ethylbenzene | ND |

Field by CLA 6/25/86

pH 11.08
Conductivity 8100

SAMPLE ID: MW - 7

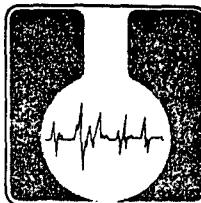
| ANALYTE | ANALYTICAL RESULTS |
|-----------------------------|--------------------|
| Methyl Bromide | ND |
| Methyl Chloride | ND |
| Methylene Chloride | ND |
| 1,1,2,2-Tetrachloroethane | ND |
| Tetrachloroethylene | ND |
| Toluene | ND |
| 1,2-Transdichloroethylene | ND |
| 1,1,1-Trichloroethane | ND |
| 1,1,2-Trichloroethane | ND |
| Trichloroethylene | ND |
| Vinyl Chloride | ND |
| Acid Compounds | |
| 2-Chlorophenol | ND |
| 2,4-Dichlorophenol | ND |
| 2,4-Dimethylphenol | ND |
| 4,6-Dinitro-o-cresol | ND |
| 2,4-Dinitrophenol | ND |
| 2-Nitrophenol | ND |
| 4-Nitrophenol | ND |
| P-chloro-m-cresol | ND |
| pentachlorophenol | ND |
| Phenol | ND |
| 2,4,6-Trichlorophenol | ND |
| Base Neutrals | |
| Acenaphthene | ND |
| Acenaphthylene | ND |
| Anthracene | ND |
| Benzidine | ND |
| Benzo(a)anthracene | 0.001 mg/l |
| Benzo(a)pyrene | ND |
| 3,4-Benzofluoranthene | ND |
| Benzo(g,h,i)perylene | ND |
| Benzo(k)fluoranthene | ND |
| Bis(2-chloroethoxy)methane | ND |
| Bis(2-chloroethyl)ether | ND |
| Bis(2-chloroisopropyl)ether | ND |
| Bis(2-ethylhexyl)phthalate | ND |
| 4-Bromophenyl phenyl ether | ND |
| Butylbenzyl phthalate | ND |
| 2-Chloronaphthalene | ND |
| 4-Chlorophenyl phenyl ether | ND |
| Chrysene | 0.002 mg/l |

SAMPLE ID: MW - 7

ANALYTE ANALYTICAL RESULTS

| | |
|----------------------------|----|
| Dibenzo(a,h)anthracene | ND |
| 1,2-Dichlorobenzene | ND |
| 1,3-Dichlorobenzene | ND |
| 1,4-Dichlorobenzene | ND |
| 3,3-Dichlorobenzidine | ND |
| Diethyl phthalate | ND |
| Dimethyl phthalate | ND |
| Din-n-butyl phthalate | ND |
| 2,4-Dinitrotoluene | ND |
| 2,6-Dinitrotoluene | ND |
| Di-n-octyl phthalate | ND |
| 1,2-Diphenylhydrazine | ND |
| Fluoranthene | ND |
| Fluorene | ND |
| Hexachlorobenzene | ND |
| Hexachlorobutadiene | ND |
| Hexachlorocyclopentadiene | ND |
| Hexachloroethane | ND |
| Indeno(1,2,3-cd)pyrene | ND |
| Isophorone | ND |
| Naphthalene | ND |
| Nitrobenzene | ND |
| N-nitrosodimethylamine | ND |
| N-nitrosodie-n-propylamine | ND |
| N-nitrosodiphenylamine | ND |
| Phenanthrene | ND |
| Pyrene | ND |
| 1,2,4-Trichlorobenzene | ND |

ND = None Detected



ASSAIGAI ANALYTICAL LABORATORIES



TO: Bloomfield Refinery
Attn: Chris Hawley
PO Box 159
Bloomfield, NM 87413

DATE: 23 July 1986
1030

SAMPLE ID: MW - 8

ANALYTE

ANALYTICAL RESULTS

| | |
|--------------------------|-------------|
| CN | <0.01 mg/l |
| TDS | 2910 mg/l |
| Cl | 839.7 mg/l |
| SO 4 | 1500 mg/l |
| Phenols | 0.005 mg/l |
| TOC | 13 mg/l |
| Sb | <0.01 mg/l |
| As | 0.072 mg/l |
| Be | <0.01 mg/l |
| Cd | <0.010 mg/l |
| Cr | <0.050 mg/l |
| Cu | <0.03 mg/l |
| Pb | 0.055 mg/l |
| Hg | <0.002 mg/l |
| Ni | 0.86 mg/l |
| Se | 0.21 mg/l |
| Ag | <0.050 mg/l |
| Tl | <0.01 mg/l |
| Zn | 0.020 mg/l |
| Volatiles | |
| Acrolein | ND |
| Acrylonitrile | ND |
| Benzene | ND |
| Bromoform | ND |
| Carbon Tetrachloride | ND |
| Chlorobenzene | ND |
| Chlorodibromomethane | ND |
| Chloroethane | ND |
| 2-Chloroethylvinyl ether | ND |
| Chloroform | ND |
| Dichlorobromomethane | ND |
| 1,1-Dichloroethane | ND |
| 1,2-Dichloroethane | ND |
| 1,1-Dichloroethylene | ND |
| 1,2-Dichloropropane | ND |
| 1,2-Dichloropropylene | ND |
| Ethylbenzene | ND |

Field by Clt 6/23/86
pH 7.26
Conductivity 4400

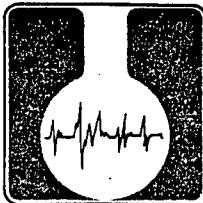
SAMPLE ID: MW - 8

| ANALYTE | ANALYTICAL RESULTS |
|-----------------------------|--------------------|
| Methyl Bromide | ND |
| Methyl Chloride | ND |
| Methylene Chloride | ND |
| 1,1,2,2-Tetrachloroethane | ND |
| Tetrachloroethylene | ND |
| Toluene | ND |
| 1,2-Transdichloroethylene | ND |
| 1,1,1-Trichloroethane | ND |
| 1,1,2-Trichloroethane | ND |
| Trichloroethylene | ND |
| Vinyl Chloride | ND |
| Acid Compounds | |
| 2-Chlorophenol | ND |
| 2,4-Dichlorophenol | ND |
| 2,4-Dimethylphenol | ND |
| 4,6-Dinitro-o-cresol | ND |
| 2,4-Dinitrophenol | ND |
| 2-Nitrophenol | ND |
| 4-Nitrophenol | ND |
| P-chloro-m-cresol | ND |
| pentachlorophenol | ND |
| Phenol | ND |
| 2,4,6-Trichlorophenol | ND |
| Base Neutrals | |
| Acenaphthene | ND |
| Acenaphthylene | ND |
| Anthracene | ND |
| Benzidine | ND |
| Benzo(a)anthracene | ND |
| Benzo(a)pyrene | ND |
| 3,4-Benzofluoranthene | ND |
| Benzo(g,h,i)perylene | ND |
| Benzo(k)fluoranthene | ND |
| Bis(2-chloroethoxy)methane | ND |
| Bis(2-chlororoethyl)ether | ND |
| Bis(2-chloroisopropyl)ether | ND |
| Bis(2-ethylhexyl)phthalate | ND |
| 4-Bromophenyl phenyl ether | ND |
| Butylbenzyl phthalate | ND |
| 2-Chloronaphthalene | ND |
| 4-Chlorophenyl phenyl ether | ND |
| Chrysene | ND |

SAMPLE ID: MW - 8

| ANALYTE | ANALYTICAL RESULTS |
|-----------------------------|--------------------|
| Dibenzo(a,h)anthracene | ND |
| 1,2-Dichlorobenzene | ND |
| 1,3-Dichlorobenzene | ND |
| 1,4-Dichlorobenzene | ND |
| 3,3-Dichlorobenzidine | ND |
| Diethyl phthalate | ND |
| Dimethyl phthalate | ND |
| Din-n-butyl phthalate | ND |
| 2,4-Dinitrotoluene | ND |
| 2,6-Dinitrotoluene | ND |
| Di-n-octyl phthalate | ND |
| 1,2-Diphenylhydrazine | ND |
| Fluoranthene | ND |
| Fluorene | ND |
| Hexachlorobenzene | ND |
| Hexachlorobutadiene | ND |
| Hexachlorocyclopentadiene | ND |
| Hexachloroethane | ND |
| Indeno(1,2,3-cd)pyrene | ND |
| Isophorone | ND |
| Naphthalene | ND |
| Nitrobenzene | ND |
| N-nitrosodimethylamine | ND |
| N-nitrosodi-e-n-propylamine | ND |
| N-nitrosodiphenylamine | ND |
| Phenanthrene | ND |
| Pyrene | ND |
| 1,2,4-Trichlorobenzene | ND |

ND = None Detected



ASSAIGAI ANALYTICAL LABORATORIES



TO: Bloomfield Refinery
Attn: Chris Hawley
PO Box 159
Bloomfield, NM 87413

DATE: 23 July 1986
1030

SAMPLE ID: MW - 9

| ANALYTE | ANALYTICAL RESULTS |
|--------------------------|--------------------|
| CN | 0.4 mg/l |
| TDS | 1718 mg/l |
| Cl | 1009.7 mg/l |
| SO ₄ | 114 mg/l |
| Phenols | 0.372 mg/l |
| TOC | 180 mg/l |
| Sb | <0.01 mg/l |
| As | <0.05 mg/l |
| Be | <0.01 mg/l |
| Cd | <0.010 mg/l |
| Cr | <0.050 mg/l |
| Cu | <0.03 mg/l |
| Pb | 0.059 mg/l |
| Hg | <0.002 mg/l |
| Ni | 0.25 mg/l |
| Se | 0.040 mg/l |
| Ag | <0.050 mg/l |
| Tl | <0.01 mg/l |
| Zn | 0.015 mg/l |
| Volatiles | |
| Acrolein | ND |
| Acrylonitrile | ND |
| Benzene | 4 mg/l |
| Bromoform | ND |
| Carbon Tetrachloride | ND |
| Chlorobenzene | ND |
| Chlorodibromomethane | ND |
| Chloroethane | ND |
| 2-Chloroethylvinyl ether | ND |
| Chloroform | ND |
| Dichlorobromomethane | ND |
| 1,1-Dichloroethane | ND |
| 1,2-Dichloroethane | ND |
| 1,1-Dichloroethylene | ND |
| 1,2-Dichloropropane | ND |
| 1,2-Dichloropropylene | ND |
| Ethylbenzene | 0.71 mg/l |

ANALYTE

ANALYTICAL RESULTS

CN 0.4 mg/l
TDS 1718 mg/l
Cl 1009.7 mg/l
SO₄ 114 mg/l
Phenols 0.372 mg/l
TOC 180 mg/l
Sb <0.01 mg/l
As <0.05 mg/l
Be <0.01 mg/l
Cd <0.010 mg/l
Cr <0.050 mg/l
Cu <0.03 mg/l
Pb 0.059 mg/l
Hg <0.002 mg/l
Ni 0.25 mg/l
Se 0.040 mg/l
Ag <0.050 mg/l
Tl <0.01 mg/l
Zn 0.015 mg/l

Volatile
Acrolein ND
Acrylonitrile ND
Benzene 4 mg/l
Bromoform ND
Carbon Tetrachloride ND
Chlorobenzene ND
Chlorodibromomethane ND
Chloroethane ND
2-Chloroethylvinyl ether ND
Chloroform ND
Dichlorobromomethane ND
1,1-Dichloroethane ND
1,2-Dichloroethane ND
1,1-Dichloroethylene ND
1,2-Dichloropropane ND
1,2-Dichloropropylene ND
Ethylbenzene 0.71 mg/l

Field by CLT 6/4/86

pH 6.98
Conductivity 2500

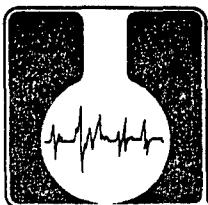
SAMPLE ID: MW - 9

| ANALYTE | ANALYTICAL RESULTS |
|-----------------------------|--------------------|
| Methyl Bromide | ND |
| Methyl Chloride | ND |
| Methylene Chloride | ND |
| 1,1,2,2-Tetrachloroethane | ND |
| Tetrachloroethylene | ND |
| Toluene | 1.7 mg/l |
| 1,2-Transdichloroethylene | ND |
| 1,1,1-Trichloroethane | ND |
| 1,1,2-Trichloroethane | ND |
| Trichloroethylene | ND |
| Vinyl Chloride | ND |
| Acid Compounds | |
| 2-Chlorophenol | ND |
| 2,4-Dichlorophenol | ND |
| 2,4-Dimethylphenol | 0.150 mg/l |
| 4,6-Dinitro-o-cresol | ND |
| 2,4-Dinitrophenol | ND |
| 2-Nitrophenol | ND |
| 4-Nitrophenol | ND |
| P-chloro-m-cresol | ND |
| pentachlorophenol | ND |
| Phenol | 0.170 mg/l |
| 2,4,6-Trichlorophenol | ND |
| Base Neutrals | |
| Acenaphthene | ND |
| Acenaphthylene | ND |
| Anthracene | ND |
| Benzidine | ND |
| Benzo(a)anthracene | ND |
| Benzo(a)pyrene | ND |
| 3,4-Benzofluoranthene | ND |
| Benzo(g,h,i)perylene | ND |
| Benzo(k)fluoranthene | ND |
| Bis(2-chloroethoxy)methane | ND |
| Bis(2-chloroethyl)ether | ND |
| Bis(2-chloroisopropyl)ether | ND |
| Bis(2-ethylhexyl)phthalate | ND |
| 4-Bromophenyl phenyl ether | ND |
| Butylbenzyl phthalate | ND |
| 2-Chloronaphthalene | ND |
| 4-Chlorophenyl phenyl ether | ND |
| Chrysene | ND |

SAMPLE ID: MW - 9

| ANALYTE | ANALYTICAL RESULTS |
|---------------------------|--------------------|
| Dibenzo(a,h)anthracene | ND |
| 1,2-Dichlorobenzene | ND |
| 1,3-Dichlorobenzene | ND |
| 1,4-Dichlorobenzene | ND |
| 3,3-Dichlorobenzidine | ND |
| Diethyl phthalate | ND |
| Dimethyl phthalate | ND |
| Din-n-butyl phthalate | ND |
| 2,4-Dinitrotoluene | ND |
| 2,6-Dinitrotoluene | ND |
| Di-n-octyl phthalate | ND |
| 1,2-Diphenylhydrazine | ND |
| Fluoranthene | ND |
| Fluorene | ND |
| Hexachlorobenzene | ND |
| Hexachlorobutadiene | ND |
| Hexachlorocyclopentadiene | ND |
| Hexachloroethane | ND |
| Indeno(1,2,3-cd)pyrene | ND |
| Isophorone | ND |
| Naphthalene | ND |
| Nitrobenzene | ND |
| N-nitrosodimethylamine | ND |
| N-nitrosodi-n-propylamine | ND |
| N-nitrosodiphenylamine | ND |
| Phenanthrene | ND |
| Pyrene | ND |
| 1,2,4-Trichlorobenzene | ND |

ND = None Detected,



ASSAIGAI ANALYTICAL LABORATORIES

TO: Bloomfield Refinery
Attn: Chris Hawley
PO Box 159
Bloomfield, NM 87413

DATE: 23 July 1986
1030

SAMPLE ID: MW - 10

| ANALYTE | ANALYTICAL RESULTS |
|--------------------------|--------------------|
| CN | <0.01 mg/l |
| TDS | 2820 mg/l |
| Cl | 569.8 mg/l |
| SO 4 | 165 mg/l |
| Phenols | 0.186 mg/l |
| TOC | 76 mg/l |
| Sb | <0.01 mg/l |
| As | 0.053 mg/l |
| Be | <0.01 mg/l |
| Cd | <0.010 mg/l |
| Cr | <0.050 mg/l |
| Cu | <0.03 mg/l |
| Pb | 0.059 mg/l |
| Hg | <0.002 mg/l |
| Ni | <0.25 mg/l |
| Se | 0.040 mg/l |
| Ag | <0.050 mg/l |
| Tl | <0.01 mg/l |
| Zn | 0.015 mg/l |
| Volatiles | |
| Acrolein | ND |
| Acrylonitrile | ND |
| Benzene | ND |
| Bromoform | ND |
| Carbon Tetrachloride | ND |
| Chlorobenzene | ND |
| Chlorodibromomethane | ND |
| Chloroethane | ND |
| 2-Chloroethylvinyl ether | ND |
| Chloroform | ND |
| Dichlorobromomethane | ND |
| 1,1-Dichloroethane | ND |
| 1,2-Dichloroethane | ND |
| 1,1-Dichloroethylene | ND |
| 1,2-Dichloropropane | ND |
| 1,2-Dichloropropylene | ND |
| Ethylbenzene | ND |

Field by C4 6/24/86

pH 7.08

Conductivity 4400

SAMPLE ID: MW - 10

| ANALYTE | ANALYTICAL RESULTS |
|-----------------------------|--------------------|
| Methyl Bromide | ND |
| Methyl Chloride | ND |
| Methylene Chloride | ND |
| 1,1,2,2-Tetrachloroethane | ND |
| Tetrachloroethylene | ND |
| Toluene | ND |
| 1,2-Transdichloroethylene | ND |
| 1,1,1-Trichloroethane | ND |
| 1,1,2-Trichloroethane | ND |
| Trichloroethylene | ND |
| Vinyl Chloride | ND |
| Acid Compounds | |
| 2-Chlorophenol | ND |
| 2,4-Dichlorophenol | ND |
| 2,4-Dimethylphenol | ND |
| 4,6-Dinitro-o-cresol | ND |
| 2,4-Dinitrophenol | ND |
| 2-Nitrophenol | ND |
| 4-Nitrophenol | ND |
| P-chloro-m-cresol | ND |
| pentachlorophenol | ND |
| Phenol | ND |
| 2,4,6-Trichlorophenol | ND |
| Base Neutrals | |
| Acenaphthene | ND |
| Acenaphthylene | ND |
| Anthracene | ND |
| Benzidine | ND |
| Benzo(a)anthracene | ND |
| Benzo(a)pyrene | ND |
| 3,4-Benzofluoranthene | ND |
| Benzo(g,h,i)perylene | ND |
| Benzo(k)fluoranthene | ND |
| Bis(2-chloroethoxy)methane | ND |
| Bis(2-chlororoethyl)ether | ND |
| Bis(2-chloroisopropyl)ether | ND |
| Bis(2-ethylhexyl)phthalate | ND |
| 4-Bromophenyl phenyl ether | ND |
| Butylbenzyl phthalate | ND |
| 2-Chloronaphthalene | ND |
| 4-Chlorophenyl phenyl ether | ND |
| Chrysene | ND |

SAMPLE ID: MW - 10

| ANALYTE | ANALYTICAL RESULTS |
|----------------------------|--------------------|
| Dibenzo(a,h)anthracene | ND |
| 1,2-Dichlorobenzene | ND |
| 1,3-Dichlorobenzene | ND |
| 1,4-Dichlorobenzene | ND |
| 3,3-Dichlorobenzidine | ND |
| Diethyl phthalate | ND |
| Dimethyl phthalate | ND |
| Din-n-butyl phthalate | ND |
| 2,4-Dinitrotoluene | ND |
| 2,6-Dinitrotoluene | ND |
| Di-n-octyl phthalate | ND |
| 1,2-Diphenylhydrazine | ND |
| Fluoranthene | ND |
| Fluorene | ND |
| Hexachlorobenzene | ND |
| Hexachlorobutadiene | ND |
| Hexachlorocyclopentadiene | ND |
| Hexachloroethane | ND |
| Indeno(1,2,3-cd)pyrene | ND |
| Isophorone | ND |
| Naphthalene | ND |
| Nitrobenzene | ND |
| N-nitrosodimethylamine | ND |
| N-nitrosodie-n-propylamine | ND |
| N-nitrosodiphenylamine | ND |
| Phenanthrene | ND |
| Pyrene | ND |
| 1,2,4-Trichlorobenzene | ND |

ND = None Detected

TO: Bloomfield Refinery
Attn: Chris Hawley
PO Box 159
Bloomfield, NM 87413

DATE: 23 July 1986
1030

NOMINAL DETECTION LIMITS

ANALYTE

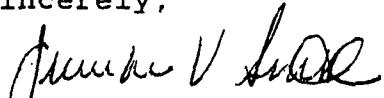
| | |
|-----------------|------------|
| CN | 0.01 mg/l |
| TDS | 1 mg/l |
| Cl | 1.0 mg/l |
| SO ₄ | 1.0 mg/l |
| Phenols | 0.002 mg/l |
| TOC | 0.1 mg/l |
| Sb | 0.01 mg/l |
| As | 0.05 mg/l |
| Be | 0.01 mg/l |
| Cd | 0.010 mg/l |
| Cr | 0.050 mg/l |
| Cu | 0.03 mg/l |
| Pb | 0.050 mg/l |
| Hg | 0.002 mg/l |
| Ni | 0.06 mg/l |
| Se | 0.010 mg/l |
| Ag | 0.050 mg/l |
| Tl | 0.01 mg/l |
| Zn | 0.01 mg/l |
| Benzene | 0.001 mg/l |
| Toluene | 0.001 mg/l |
| Xylenes | 0.001 mg/l |
| Ethylbenzene | 0.001 mg/l |
| Ba | 0.01 mg/l |
| Fe | 0.04 mg/l |
| Mn | 0.005 mg/l |
| Al | 0.05 mg/l |
| B | 0.01 mg/l |
| Co | 0.05 mg/l |
| Mo | 0.01 mg/l |
| F | 0.1 mg/l |
| No 3 as N | 0.01 mg/l |
| 1,2-DCE | 0.001 mg/l |
| 1,1-DCE | 0.001 mg/l |
| 1,1,2,2-TCE | 0.001 mg/l |
| 1,1,2-TCE | 0.001 mg/l |

Detection limits for Volatiles, Base/Neutrals and Acid Compounds all 0.001 mg/l

REFERENCE: "Test Methods for Evaluating Solid Waste,
Physical/Chemical Methods", USEPA, SW 846, EMSL-Cincinnati,
1982.

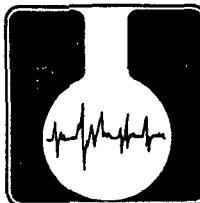
An invoice for services is enclosed. Thank you for contacting
Assaigai Laboratories.

Sincerely,



Jennifer V. Smith, Ph.D.
Laboratory Director

THIRD QUARTER



ASSAIGAI ANALYTICAL LABORATORIES

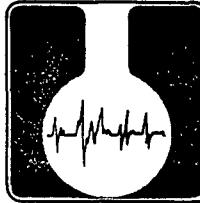


TO: Bloomfield Refinery
Attn: Chris Hawley
PO Box 159
Bloomfield NM 87413

DATE: 11 November 1986
1624

SAMPLE ID: MW - 1

| ANALYTE | ANALYTICAL RESULTS |
|-----------------|--------------------|
| CN | 0.07 mg/l |
| Phenols | 0.19 mg/l |
| TOC | 24 mg/l |
| TDS | 2866 mg/l |
| Cl | 814 mg/l |
| SO ₄ | 673 mg/l |
| Sb | <0.01 mg/l |
| As | 0.05 mg/l |
| Be | <0.01 mg/l |
| Cd | <0.01 mg/l |
| Cr | <0.05 mg/l |
| Cu | <0.03 mg/l |
| Pb | 0.15 mg/l |
| Hg | <0.002 mg/l |
| Ni | 0.07 mg/l |
| Se | 0.033 mg/l |
| Ag | <0.05 mg/l |
| Tl | <0.01 mg/l |
| Zn | 0.04 mg/l |
| Benzene | <0.001 mg/l |
| Toluene | <0.001 mg/l |
| Ethylbenzene | <0.001 mg/l |
| Xylenes | <0.001 mg/l |



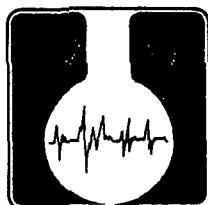
ASSAIGAI ANALYTICAL LABORATORIES

TO: Bloomfield Refinery
Attn: Chris Hawley
PO Box 159
Bloomfield NM 87413

DATE: 11 November 1986
1624

SAMPLE ID: MW - 2

| ANALYTE | ANALYTICAL RESULTS |
|-----------------|--------------------|
| CN | 0.18 mg/l |
| Phenols | 0.17 mg/l |
| TOC | 23 mg/l |
| TDS | 3598 mg/l |
| Cl | 993 mg/l |
| SO ₄ | 1104 mg/l |
| Sb | <0.01 mg/l |
| As | 0.08 mg/l |
| Be | <0.01 mg/l |
| Cd | 0.03 mg/l |
| Cr | <0.05 mg/l |
| Cu | <0.03 mg/l |
| Pb | 0.08 mg/l |
| Hg | <0.002 mg/l |
| Ni | 0.12 mg/l |
| Se | 0.104 mg/l |
| Ag | <0.05 mg/l |
| Tl | <0.01 mg/l |
| Zn | 0.02 mg/l |
| Benzene | <0.001 mg/l |
| Toluene | <0.001 mg/l |
| Ethylbenzene | <0.001 mg/l |
| Xylenes | <0.001 mg/l |



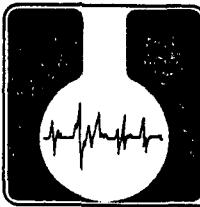
ASSAIGAI ANALYTICAL LABORATORIES

TO: Bloomfield Refinery
Attn: Chris Hawley
PO Box 159
Bloomfield NM 87413

DATE: 11 November 1986
1624

SAMPLE ID: MW - 3

| ANALYTE | ANALYTICAL RESULTS |
|-----------------|--------------------|
| CN | 0.17 mg/l |
| Phenols | 0.082 mg/l |
| TOC | 16 mg/l |
| TDS | 5514 mg/l |
| Cl | 1290 mg/l |
| SO ₄ | 2056 mg/l |
| Sb | <0.01 mg/l |
| As | 0.21 mg/l |
| Be | <0.01 mg/l |
| Cd | <0.01 mg/l |
| Cr | <0.05 mg/l |
| Cu | <0.03 mg/l |
| Pb | 0.18 mg/l |
| Hg | <0.002 mg/l |
| Ni | 0.14 mg/l |
| Se | 0.100 mg/l |
| Ag | <0.05 mg/l |
| Tl | <0.01 mg/l |
| Zn | 0.02 mg/l |
| Benzene | <0.001 mg/l |
| Toluene | <0.001 mg/l |
| Ethylbenzene | <0.001 mg/l |
| Xylenes | <0.001 mg/l |



ASSAIGAI ANALYTICAL LABORATORIES

TO: Bloomfield Refinery
Attn: Chris Hawley
PO Box 159
Bloomfield NM 87413

DATE: 11 November 1986
1624

SAMPLE ID: MW - 4

ANALYTE

ANALYTICAL RESULTS

| | |
|--------------------------|-------------|
| CN | <0.01 mg/l |
| Phenols | 0.085 mg/l |
| TOC | 63 mg/l |
| TDS | 2308 mg/l |
| Cl | 754 mg/l |
| SO ₄ | <0.01 mg/l |
| Sb | <0.01 mg/l |
| As | 0.08 mg/l |
| Be | <0.01 mg/l |
| Cd | <0.01 mg/l |
| Cr | <0.05 mg/l |
| Cu | <0.03 mg/l |
| Pb | <0.05 mg/l |
| Hg | <0.002 mg/l |
| Ni | 0.12 mg/l |
| Se | 0.063 mg/l |
| Ag | <0.05 mg/l |
| Tl | <0.01 mg/l |
| Zn | <0.008 mg/l |
| Volatiles | |
| Acrolein | ND |
| Acrylonitrile | ND |
| Benzene | 6.65 mg/l |
| Bromoform | ND |
| Carbon Tetrachloride | ND |
| Chlorobenzene | ND |
| Chlorodibromomethane | ND |
| Chloroethane | ND |
| 2-Chloroethylvinyl ether | ND |
| Chloroform | ND |
| Dichlorobromomethane | ND |
| 1,1-Dichloroethane | ND |
| 1,2-Dichloroethane | ND |
| 1,1-Dichloroethylene | ND |
| 1,2-Dichloropropane | ND |
| 1,2-Dichloropropylene | ND |
| Ethylbenzene | 0.140 mg/l |

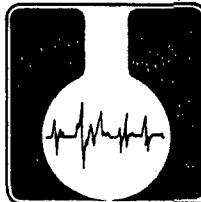
SAMPLE ID: MW - 4

| ANALYTE | ANALYTICAL RESULTS |
|-----------------------------|--------------------|
| Methyl Bromide | ND |
| Methyl Chloride | ND |
| Methylene Chloride | ND |
| 1,1,2,2-Tetrachloroethane | ND |
| Tetrachloroethylene | ND |
| Toluene | 0.407 mg/l |
| 1,2-Transdichloroethylene | ND |
| 1,1,1-Trichloroethane | ND |
| 1,1,2-Trichloroethane | ND |
| Trichloroethylene | ND |
| Vinyl Chloride | ND |
| Acid Compounds | |
| 2-Chlorophenol | 0.001 mg/l |
| 2,4-Dichlorophenol | ND |
| 2,4-Dimethylphenol | ND |
| 4,6-Dinitro-o-cresol | ND |
| 2,4-Dinitrophenol | ND |
| 2-Nitrophenol | 0.026 mg/l |
| 4-Nitrophenol | 0.331 mg/l |
| P-chloro-m-cresol | 0.045 mg/l |
| pentachlorophenol | ND |
| Phenol | ND |
| 2,4,6-Trichlorophenol | ND |
| Base Neutrals | |
| Acenaphthene | ND |
| Acenaphthylene | ND |
| Anthracene | ND |
| Benzidine | ND |
| Benzo(a)anthracene | 0.010 mg/l |
| Benzo(a)pyrene | ND |
| 3,4-Benzofluoranthene | ND |
| Benzo(g,h,i)perylene | ND |
| Benzo(k)fluoranthene | ND |
| Bis(2-chloroethoxy)methane | ND |
| Bis(2-chloroethyl)ether | ND |
| Bis(2-chloroisopropyl)ether | ND |
| Bis(2-ethylhexyl)phthalate | ND |
| 4-Bromophenyl phenyl ether | ND |
| Butylbenzyl phthalate | ND |
| 2-Chloronaphthalene | ND |
| 4-Chlorophenyl phenyl ether | ND |
| Chrysene | ND |

SAMPLE ID: MW - 4

| ANALYTE | ANALYTICAL RESULTS |
|----------------------------|--------------------|
| Dibenzo(a,h)anthracene | ND |
| 1,2-Dichlorobenzene | ND |
| 1,3-Dichlorobenzene | ND |
| 1,4-Dichlorobenzene | ND |
| 3,3-Dichlorobenzidine | ND |
| Diethyl phthalate | ND |
| Dimethyl phthalate | ND |
| Di-n-butyl phthalate | ND |
| 2,4-Dinitrotoluene | ND |
| 2,6-Dinitrotoluene | ND |
| Di-n-octyl phthalate | ND |
| 1,2-Diphenylhydrazine | ND |
| Fluoranthene | ND |
| Fluorene | ND |
| Hexachlorobenzene | ND |
| Hexachlorobutadiene | ND |
| Hexachlorocyclopentadiene | ND |
| Hexachloroethane | ND |
| Indeno(1,2,3-cd)pyrene | ND |
| Isophorone | ND |
| Naphthalene | 0.015 mg/l |
| Nitrobenzene | ND |
| N-nitrosodimethylamine | ND |
| N-nitrosodie-n-propylamine | ND |
| N-nitrosodiphenylamine | ND |
| Phenanthrene | ND |
| Pyrene | 0.005 mg/l |
| 1,2,4-Trichlorobenzene | ND |

ND = None Detected



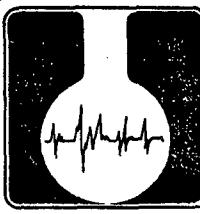
ASSAIGAI ANALYTICAL LABORATORIES

TO: Bloomfield Refinery
Attn: Chris Hawley
PO Box 159
Bloomfield NM 87413

DATE: 11 November 1986
1624

SAMPLE ID: MW - 5

| ANALYTE | ANALYTICAL RESULTS |
|-----------------|--------------------|
| CN | 0.24 mg/l |
| Phenols | 0.034 mg/l |
| TOC | 20 mg/l |
| TDS | 3184 mg/l |
| Cl | 1151 mg/l |
| SO ₄ | 1237 mg/l |
| Sb | <0.01 mg/l |
| As | 0.07 mg/l |
| Be | <0.01 mg/l |
| Cd | <0.01 mg/l |
| Cr | <0.05 mg/l |
| Cu | <0.03 mg/l |
| Pb | <0.05 mg/l |
| Hg | <0.002 mg/l |
| Ni | 0.09 mg/l |
| Se | 0.030 mg/l |
| Ag | <0.05 mg/l |
| Tl | <0.01 mg/l |
| Zn | 0.02 mg/l |
| Benzene | <0.001 mg/l |
| Toluene | <0.001 mg/l |
| Ethylbenzene | <0.001 mg/l |
| Xylenes | <0.001 mg/l |



ASSAIGAI ANALYTICAL LABORATORIES



TO: Bloomfield Refinery
Attn: Chris Hawley
PO Box 159
Bloomfield NM 87413

DATE: 11 November 1986
1624

SAMPLE ID: MW - 7

ANALYTE

ANALYTICAL RESULTS

| | |
|--------------------------|-------------|
| CN | 0.10 mg/l |
| Phenols | 0.036 mg/l |
| TOC | 4 mg/l |
| TDS | 6348 mg/l |
| Cl | 20 mg/l |
| SO ₄ | 5802 mg/l |
| Sb | <0.01 mg/l |
| As | 0.22 mg/l |
| Be | <0.01 mg/l |
| Cd | <0.01 mg/l |
| Cr | <0.05 mg/l |
| Cu | <0.03 mg/l |
| Pb | 0.05 mg/l |
| Hg | <0.002 mg/l |
| Ni | 0.08 mg/l |
| Se | 0.36 mg/l |
| Ag | <0.05 mg/l |
| Tl | <0.01 mg/l |
| Zn | 0.02 mg/l |
| Volatiles | |
| Acrolein | ND |
| Acrylonitrile | ND |
| Benzene | 0.058 mg/l |
| Bromoform | ND |
| Carbon Tetrachloride | ND |
| Chlorobenzene | ND |
| Chlorodibromomethane | ND |
| Chloroethane | ND |
| 2-Chloroethylvinyl ether | ND |
| Chloroform | ND |
| Dichlorobromomethane | ND |
| 1,1-Dichloroethane | ND |
| 1,2-Dichloroethane | ND |
| 1,1-Dichloroethylene | ND |
| 1,2-Dichloropropane | ND |
| 1,2-Dichloropropylene | ND |
| Ethylbenzene | 0.004 mg/l |

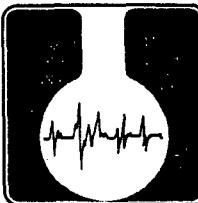
SAMPLE ID: MW -7

| ANALYTE | ANALYTICAL RESULTS |
|-----------------------------|--------------------|
| Methyl Bromide | ND |
| Methyl Chloride | ND |
| Methylene Chloride | ND |
| 1,1,2,2-Tetrachloroethane | ND |
| Tetrachloroethylene | ND |
| Toluene | 0.006 mg/l |
| 1,2-Transdichloroethylene | ND |
| 1,1,1-Trichloroethane | ND |
| 1,1,2-Trichloroethane | ND |
| Trichloroethylene | ND |
| Vinyl Chloride | ND |
| Acid Compounds | |
| 2-Chlorophenol | ND |
| 2,4-Dichlorophenol | ND |
| 2,4-Dimethylphenol | ND |
| 4,6-Dinitro-o-cresol | ND |
| 2,4-Dinitrophenol | ND |
| 2-Nitrophenol | ND |
| 4-Nitrophenol | 0.007 mg/l |
| P-chloro-m-cresol | 0.001 mg/l |
| pentachlorophenol | ND |
| Phenol | ND |
| 2,4,6-Trichlorophenol | ND |
| Base Neutrals | |
| Acenaphthene | ND |
| Acenaphthylene | ND |
| Anthracene | ND |
| Benzidine | ND |
| Benzo(a)anthracene | ND |
| Benzo(a)pyrene | ND |
| 3,4-Benzofluoranthene | ND |
| Benzo(g,h,i)perylene | ND |
| Benzo(k)fluoranthene | ND |
| Bis(2-chloroethoxy)methane | ND |
| Bis(2-chloroethyl)ether | ND |
| Bis(2-chloroisopropyl)ether | ND |
| Bis(2-ethylhexyl)phthalate | ND |
| 4-Bromophenyl phenyl ether | ND |
| Butylbenzyl phthalate | ND |
| 2-Chloronaphthalene | ND |
| 4-Chlorophenyl phenyl ether | ND |
| Chrysene | ND |

SAMPLE ID: MW - 7

| ANALYTE | ANALYTICAL RESULTS |
|----------------------------|--------------------|
| Dibenzo(a,h)anthracene | ND |
| 1,2-Dichlorobenzene | ND |
| 1,3-Dichlorobenzene | ND |
| 1,4-Dichlorobenzene | ND |
| 3,3-Dichlorobenzidine | ND |
| Diethyl phthalate | ND |
| Dimethyl phthalate | ND |
| Di-n-butyl phthalate | ND |
| 2,4-Dinitrotoluene | ND |
| 2,6-Dinitrotoluene | ND |
| Di-n-octyl phthalate | ND |
| 1,2-Diphenylhydrazine | ND |
| Fluoranthene | ND |
| Fluorene | ND |
| Hexachlorobenzene | ND |
| Hexachlorobutadiene | ND |
| Hexachlorocyclopentadiene | ND |
| Hexachloroethane | ND |
| Indeno(1,2,3-cd)pyrene | ND |
| Isophorone | ND |
| Naphthalene | ND |
| Nitrobenzene | ND |
| N-nitrosodimethylamine | ND |
| N-nitrosodie-n-propylamine | ND |
| N-nitrosodiphenylamine | ND |
| Phenanthrene | ND |
| Pyrene | ND |
| 1,2,4-Trichlorobenzene | ND |

ND = None Detected



ASSAIGAI ANALYTICAL LABORATORIES

TO: Bloomfield Refinery
Attn: Chris Hawley
PO Box 159
Bloomfield NM 87413

DATE: 11 November 1986
1624

SAMPLE ID: MW - 8

| ANALYTE | ANALYTICAL RESULTS |
|--------------------------|--------------------|
| CN | <0.01 mg/l |
| Phenols | 0.097 mg/l |
| TOC | 8 mg/l |
| TDS | 2284 mg/l |
| Cl | 576 mg/l |
| SO 4 | 586 mg/l |
| Sb | <0.01 mg/l |
| As | 0.03 mg/l |
| Be | <0.01 mg/l |
| Cd | <0.01 mg/l |
| Cr | <0.05 mg/l |
| Cu | <0.03 mg/l |
| Pb | <0.05 mg/l |
| Hg | <0.002 mg/l |
| Ni | 0.21 mg/l |
| Se | <0.01 mg/l |
| Ag | <0.05 mg/l |
| Tl | <0.01 mg/l |
| Zn | 0.02 mg/l |
| Volatiles | |
| Acrolein | ND |
| Acrylonitrile | ND |
| Benzene | ND |
| Bromoform | ND |
| Carbon Tetrachloride | ND |
| Chlorobenzene | ND |
| Chlorodibromomethane | ND |
| Chloroethane | ND |
| 2-Chloroethylvinyl ether | ND |
| Chloroform | ND |
| Dichlorobromomethane | ND |
| 1,1-Dichloroethane | ND |
| 1,2-Dichloroethane | ND |
| 1,1-Dichloroethylene | ND |
| 1,2-Dichloropropane | ND |
| 1,2-Dichloropropylene | ND |
| Ethylbenzene | ND |

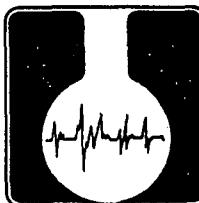
SAMPLE ID: MW - 8

| ANALYTE | ANALYTICAL RESULTS |
|-----------------------------|--------------------|
| Methyl Bromide | ND |
| Methyl Chloride | ND |
| Methylene Chloride | ND |
| 1,1,2,2-Tetrachloroethane | ND |
| Tetrachloroethylene | ND |
| Toluene | ND |
| 1,2-Transdichloroethylene | ND |
| 1,1,1-Trichloroethane | ND |
| 1,1,2-Trichloroethane | ND |
| Trichloroethylene | ND |
| Vinyl Chloride | ND |
| Acid Compounds | |
| 2-Chlorophenol | ND |
| 2,4-Dichlorophenol | ND |
| 2,4-Dimethylphenol | ND |
| 4,6-Dinitro-o-cresol | ND |
| 2,4-Dinitrophenol | ND |
| 2-Nitrophenol | ND |
| 4-Nitrophenol | 0.008 mg/l |
| P-chloro-m-cresol | ND |
| pentachlorophenol | ND |
| Phenol | ND |
| 2,4,6-Trichlorophenol | ND |
| Base Neutrals | |
| Acenaphthene | ND |
| Acenaphthylene | ND |
| Anthracene | ND |
| Benzidine | ND |
| Benzo(a)anthracene | ND |
| Benzo(a)pyrene | ND |
| 3,4-Benzofluoranthene | ND |
| Benzo(g,h,i)perylene | ND |
| Benzo(k)fluoranthene | ND |
| Bis(2-chloroethoxy)methane | ND |
| Bis(2-chloroethyl)ether | ND |
| Bis(2-chloroisopropyl)ether | ND |
| Bis(2-ethylhexyl)phthalate | ND |
| 4-Bromophenyl phenyl ether | ND |
| Butylbenzyl phthalate | ND |
| 2-Chloronaphthalene | ND |
| 4-Chlorophenyl phenyl ether | ND |
| Chrysene | ND |

SAMPLE ID: MW - 8

| ANALYTE | ANALYTICAL RESULTS |
|----------------------------|--------------------|
| Dibenzo(a,h)anthracene | ND |
| 1,2-Dichlorobenzene | ND |
| 1,3-Dichlorobenzene | ND |
| 1,4-Dichlorobenzene | ND |
| 3,3-Dichlorobenzidine | ND |
| Diethyl phthalate | ND |
| Dimethyl phthalate | ND |
| Di-n-butyl phthalate | ND |
| 2,4-Dinitrotoluene | ND |
| 2,6-Dinitrotoluene | ND |
| Di-n-octyl phthalate | ND |
| 1,2-Diphenylhydrazine | ND |
| Fluoranthene | ND |
| Fluorene | ND |
| Hexachlorobenzene | ND |
| Hexachlorobutadiene | ND |
| Hexachlorocyclopentadiene | ND |
| Hexachloroethane | ND |
| Indeno(1,2,3-cd)pyrene | ND |
| Isophorone | ND |
| Naphthalene | ND |
| Nitrobenzene | ND |
| N-nitrosodimethylamine | ND |
| N-nitrosodie-n-propylamine | ND |
| N-nitrosodiphenylamine | ND |
| Phenanthrene | ND |
| Pyrene | ND |
| 1,2,4-Trichlorobenzene | ND |

ND = None Detected



ASSAIGAI ANALYTICAL LABORATORIES

TO: Bloomfield Refinery
Attn: Chris Hawley
PO Box 159
Bloomfield NM 87413

DATE: 11 November 1986
1624

SAMPLE ID: MW - 9

| ANALYTE | ANALYTICAL RESULTS |
|--------------------------|--------------------|
| CN | <0.01 mg/l |
| Phenols | 0.17 mg/l |
| TOC | 240 mg/l |
| TDS | 1428 mg/l |
| Cl | 89 mg/l |
| SO ₄ | <0.01 mg/l |
| Sb | <0.01 mg/l |
| As | 0.02 mg/l |
| Be | <0.01 mg/l |
| Cd | <0.01 mg/l |
| Cr | <0.05 mg/l |
| Cu | <0.03 mg/l |
| Pb | <0.05 mg/l |
| Hg | <0.002 mg/l |
| Ni | 0.13 mg/l |
| Se | <0.01 mg/l |
| Ag | <0.05 mg/l |
| Tl | <0.01 mg/l |
| Zn | 0.05 mg/l |
| Volatiles | |
| Acrolein | ND |
| Acrylonitrile | ND |
| Benzene | 17.7 mg/l |
| Bromoform | ND |
| Carbon Tetrachloride | ND |
| Chlorobenzene | ND |
| Chlorodibromomethane | ND |
| Chloroethane | ND |
| 2-Chloroethylvinyl ether | ND |
| Chloroform | ND |
| Dichlorobromomethane | ND |
| 1,1-Dichloroethane | ND |
| 1,2-Dichloroethane | ND |
| 1,1-Dichloroethylene | ND |
| 1,2-Dichloropropane | ND |
| 1,2-Dichloropropylene | ND |
| Ethylbenzene | 0.015 mg/l |

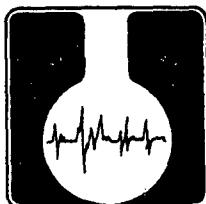
SAMPLE ID: MW - 9

| ANALYTE | ANALYTICAL RESULTS |
|-----------------------------|--------------------|
| Methyl Bromide | ND |
| Methyl Chloride | ND |
| Methylene Chloride | ND |
| 1,1,2,2-Tetrachloroethane | ND |
| Tetrachloroethylene | ND |
| Toluene | 10.6 mg/l |
| 1,2-Transdichloroethylene | ND |
| 1,1,1-Trichloroethane | ND |
| 1,1,2-Trichloroethane | ND |
| Trichloroethylene | ND |
| Vinyl Chloride | ND |
| Acid Compounds | |
| 2-Chlorophenol | ND |
| 2,4-Dichlorophenol | ND |
| 2,4-Dimethylphenol | ND |
| 4,6-Dinitro-o-cresol | ND |
| 2,4-Dinitrophenol | ND |
| 2-Nitrophenol | ND |
| 4-Nitrophenol | 1.10 mg/l |
| P-chloro-m-cresol | ND |
| pentachlorophenol | ND |
| Phenol | 0.013 mg/l |
| 2,4,6-Trichlorophenol | ND |
| Base Neutrals | |
| Acenaphthene | ND |
| Acenaphthylene | 0.028 mg/l |
| Anthracene | ND |
| Benzidine | ND |
| Benzo(a)anthracene | 0.007 mg/l |
| Benzo(a)pyrene | ND |
| 3,4-Benzofluoranthene | ND |
| Benzo(g,h,i)perylene | ND |
| Benzo(k)fluoranthene | ND |
| Bis(2-chloroethoxy)methane | ND |
| Bis(2-chloroethyl)ether | ND |
| Bis(2-chloroisopropyl)ether | ND |
| Bis(2-ethylhexyl)phthalate | ND |
| 4-Bromophenyl phenyl ether | ND |
| Butylbenzyl phthalate | ND |
| 2-Chloronaphthalene | ND |
| 4-Chlorophenyl phenyl ether | ND |
| Chrysene | ND |

SAMPLE ID: MW - 9

| ANALYTE | ANALYTICAL RESULTS |
|---------------------------|--------------------|
| Dibenzo(a,h)anthracene | ND |
| 1,2-Dichlorobenzene | ND |
| 1,3-Dichlorobenzene | ND |
| 1,4-Dichlorobenzene | ND |
| 3,3-Dichlorobenzidine | ND |
| Diethyl phthalate | ND |
| Dimethyl phthalate | ND |
| Di-n-butyl phthalate | ND |
| 2,4-Dinitrotoluene | ND |
| 2,6-Dinitrotoluene | ND |
| Di-n-octyl phthalate | ND |
| 1,2-Diphenylhydrazine | ND |
| Fluoranthene | ND |
| Fluorene | ND |
| Hexachlorobenzene | ND |
| Hexachlorobutadiene | ND |
| Hexachlorocyclopentadiene | ND |
| Hexachloroethane | ND |
| Indeno(1,2,3-cd)pyrene | ND |
| Isophorone | ND |
| Naphthalene | ND |
| Nitrobenzene | ND |
| N-nitrosodimethylamine | ND |
| N-nitrosodi-n-propylamine | ND |
| N-nitrosodiphenylamine | ND |
| Phenanthrene | ND |
| Pyrene | 0.010 mg/l |
| 1,2,4-Trichlorobenzene | ND |

ND = None Detected



ASSAIGAI ANALYTICAL LABORATORIES

TO: Bloomfield Refinery
Attn: Chris Hawley
PO Box 159
Bloomfield NM 87413

DATE: 11 November 1986
1624

SAMPLE ID: MW - 10

| ANALYTE | ANALYTICAL RESULTS |
|--------------------------|--------------------|
| CN | 0.050 mg/l |
| Phenols | 0.065 mg/l |
| TOC | 125 mg/l |
| TDS | 2408 mg/l |
| Cl | 587 mg/l |
| SO ₄ | <0.01 mg/l |
| Sb | <0.01 mg/l |
| As | 0.05 mg/l |
| Be | <0.01 mg/l |
| Cd | <0.01 mg/l |
| Cr | <0.05 mg/l |
| Cu | <0.03 mg/l |
| Pb | 0.05 mg/l |
| Hg | <0.002 mg/l |
| Ni | 0.18 mg/l |
| Se | 0.071 mg/l |
| Ag | <0.05 mg/l |
| Tl | <0.01 mg/l |
| Zn | 0.16 mg/l |
| Volatiles | |
| Acrolein | ND |
| Acrylonitrile | ND |
| Benzene | 0.041 mg/l |
| Bromoform | ND |
| Carbon Tetrachloride | ND |
| Chlorobenzene | ND |
| Chlorodibromomethane | ND |
| Chloroethane | ND |
| 2-Chloroethylvinyl ether | ND |
| Chloroform | ND |
| Dichlorobromomethane | ND |
| 1,1-Dichloroethane | ND |
| 1,2-Dichloroethane | ND |
| 1,1-Dichloroethylene | ND |
| 1,2-Dichloropropane | ND |
| 1,2-Dichloropropylene | ND |
| Ethylbenzene | ND |

SAMPLE ID: MW - 10

| ANALYTE | ANALYTICAL RESULTS |
|-----------------------------|--------------------|
| Methyl Bromide | ND |
| Methyl Chloride | ND |
| Methylene Chloride | ND |
| 1,1,2,2-Tetrachloroethane | ND |
| Tetrachloroethylene | ND |
| Toluene | 0.054 mg/l |
| 1,2-Transdichloroethylene | ND |
| 1,1,1-Trichloroethane | ND |
| 1,1,2-Trichloroethane | ND |
| Trichloroethylene | ND |
| Vinyl Chloride | ND |
| Acid Compounds | |
| 2-Chlorophenol | ND |
| 2,4-Dichlorophenol | ND |
| 2,4-Dimethylphenol | ND |
| 4,6-Dinitro-o-cresol | ND |
| 2,4-Dinitrophenol | ND |
| 2-Nitrophenol | 0.002 mg/l |
| 4-Nitrophenol | 0.016 mg/l |
| P-chloro-m-cresol | ND |
| pentachlorophenol | ND |
| Phenol | ND |
| 2,4,6-Trichlorophenol | ND |
| Base Neutrals | |
| Acenaphthene | ND |
| Acenaphthylene | ND |
| Anthracene | ND |
| Benzidine | ND |
| Benzo(a)anthracene | ND |
| Benzo(a)pyrene | ND |
| 3,4-Benzofluoranthene | ND |
| Benzo(g,h,i)perylene | ND |
| Benzo(k)fluoranthene | ND |
| Bis(2-chloroethoxy)methane | ND |
| Bis(2-chloroethyl)ether | ND |
| Bis(2-chloroisopropyl)ether | ND |
| Bis(2-ethylhexyl)phthalate | ND |
| 4-Bromophenyl phenyl ether | ND |
| Butylbenzyl phthalate | ND |
| 2-Chloronaphthalene | ND |
| 4-Chlorophenyl phenyl ether | ND |
| Chrysene | ND |

SAMPLE ID: MW - 10

| ANALYTE | ANALYTICAL RESULTS |
|----------------------------|--------------------|
| Dibenzo(a,h)anthracene | ND |
| 1,2-Dichlorobenzene | ND |
| 1,3-Dichlorobenzene | ND |
| 1,4-Dichlorobenzene | ND |
| 3,3-Dichlorobenzidine | ND |
| Diethyl phthalate | ND |
| Dimethyl phthalate | ND |
| Di-n-butyl phthalate | ND |
| 2,4-Dinitrotoluene | ND |
| 2,6-Dinitrotoluene | ND |
| Di-n-octyl phthalate | ND |
| 1,2-Diphenylhydrazine | ND |
| Fluoranthene | ND |
| Fluorene | ND |
| Hexachlorobenzene | ND |
| Hexachlorobutadiene | ND |
| Hexachlorocyclopentadiene | ND |
| Hexachloroethane | ND |
| Indeno(1,2,3-cd)pyrene | ND |
| Isophorone | ND |
| Naphthalene | ND |
| Nitrobenzene | ND |
| N-nitrosodimethylamine | ND |
| N-nitrosodie-n-propylamine | ND |
| N-nitrosodiphenylamine | ND |
| Phenanthrene | ND |
| Pyrene | ND |
| 1,2,4-Trichlorobenzene | ND |

ND = None Detected

TO: Bloomfield Refinery
Attn: Chris Hawley
PO Box 159
Bloomfield NM 87413

DATE: 11 November 1986
1624

SAMPLE ID: NOMINAL DETECTION LIMITS

ANALYTE ANALYTICAL RESULTS

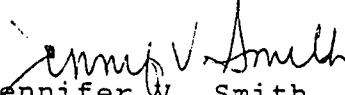
| | |
|-----------------|------------|
| CN | 0.01 mg/l |
| Phenols | 0.001 mg/l |
| TOC | 0.1 mg/l |
| TDS | 1 mg/l |
| Cl | 1.0 mg/l |
| SO ₄ | 0.01 mg/l |
| Sb | 0.01 mg/l |
| As | 0.002 mg/l |
| Be | 0.01 mg/l |
| Cd | 0.01 mg/l |
| Cr | 0.05 mg/l |
| Cu | 0.03 mg/l |
| Pb | 0.05 mg/l |
| Hg | 0.002 mg/l |
| Ni | 0.01 mg/l |
| Se | 0.01 mg/l |
| Ag | 0.05 mg/l |
| Tl | 0.01 mg/l |
| Zn | 0.008 mg/l |

Detection limits for Volatiles, Base/Neutrals and Acid Compounds all 0.001 mg/l.

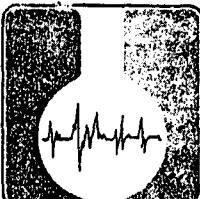
REFERENCE: "Test Methods for Evaluating Solid Waste, Physical/Chemical Method", USEPA, SW 846, EMSL-Cincinnati, 1982.

An invoice for services is enclosed. Thank you for contacting Assaigai Laboratories.

Sincerely,


Jennifer V. Smith, Ph.D.
Laboratory Director

FOURTH QUARTER



ASSAIGAI ANALYTICAL LABORATORIES



TO: Bloomfield Refinery
ATTN: Chris Hawley
PO Box 159
Bloomfield, NM 87413

DATE: 19 January 1987
2160

SAMPLE ID: MW-1

| ANALYTE | ANALYTICAL RESULTS | NOMINAL DETECTION LIMITS |
|----------------------|--------------------|--------------------------|
| CN | <0.01 mg/l | 0.01 mg/l |
| Phenols | 0.012 mg/l | 0.001 mg/l |
| TOC | 18 mg/l | 0.1 mg/l |
| TDS | 2498 mg/l | 1 mg/l |
| Cl | 774 mg/l | 1.0 mg/l |
| SO 4 | 579 mg/l | 1.0 mg/l |
| Benzene | <0.001 mg/l | 0.001 mg/l |
| Toluene | <0.001 mg/l | 0.001 mg/l |
| Xylenes | <0.001 mg/l | 0.001 mg/l |
| Ethyl Benzene | <0.001 mg/l | 0.001 mg/l |
| Sb | 0.25 mg/l | 0.002 mg/l |
| As | <0.05 mg/l | 0.05 mg/l |
| Be | 0.02 mg/l | 0.01 mg/l |
| Cd | <0.01 mg/l | 0.01 mg/l |
| Cr | <0.05 mg/l | 0.05 mg/l |
| Cu | <0.03 mg/l | 0.03 mg/l |
| Pb | <0.05 mg/l | 0.05 mg/l |
| Hg | <0.002 mg/l | 0.002 mg/l |
| Ni | 0.06 mg/l | 0.06 mg/l |
| Se | 0.03 mg/l | 0.002 mg/l |
| Ag | <0.05 mg/l | 0.05 mg/l |
| Tl | <0.1 mg/l | 0.1 mg/l |
| Zn | 0.012 mg/l | 0.001 mg/l |
| Ba | 0.055 mg/l | 0.005 mg/l |
| Fe | <0.3 mg/l | 0.3 mg/l |
| Mn | 1.11 mg/l | 0.005 mg/l |
| Al | 4.54 mg/l | 0.05 mg/l |
| B | 0.27 mg/l | 0.004 mg/l |
| Co | <0.05 mg/l | 0.05 mg/l |
| Mo | 0.17 mg/l | 0.01 mg/l |
| NO ₃ as N | 2.9 mg/l | 0.1 mg/l |
| F | 0.960 mg/l | 0.01 mg/l |
| 1,2-dichloroethane | 0.002 mg/l | 0.001 mg/l |
| 1,1-dichloroethylene | <0.001 mg/l | 0.001 mg/l |
| Tetrachloroethylene | <0.001 mg/l | 0.001 mg/l |
| Trichloroethylene | <0.001 mg/l | 0.001 mg/l |
| CCl 4 | <0.001 mg/l | 0.001 mg/l |

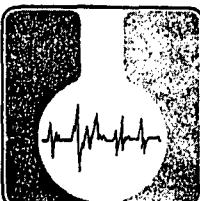
7300 Jefferson, N.E. • Albuquerque, New Mexico 87109 • (505) 345-8964

REFERENCES: "Test Methods for Evaluatin Solid Waste,-Physical/Chemical Methods", USEPA, SW 846, EMSL-Cincinnati, 1982.

An invoice for services is enclosed. Thank you for contacting Assaigai Laboratories.

Sincerely,

Jennifer V. Smith
Jennifer V. Smith, Ph.D.
Laboratory Director



ASSAIGAI ANALYTICAL LABORATORIES



TO: Bloomfield Refinery
ATTN: Chris Hawley
PO Box 159
Bloomfield, NM 87413

DATE: 19 January 1987
2160

SAMPLE ID: MW-2

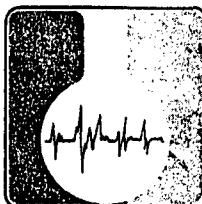
| ANALYTE | ANALYTICAL RESULTS | NOMINAL DETECTION LIMITS |
|---------------|--------------------|--------------------------|
| CN | <0.01 mg/l | 0.01 mg/l |
| Phenols | 0.110 mg/l | 0.001 mg/l |
| TOC | 15 mg/l | 0.1 mg/l |
| TDS | 3644 mg/l | 1 mg/l |
| Cl | 1012 mg/l | 1.0 mg/l |
| SO 4 | 1372 mg/l | 1.0 mg/l |
| Benzene | <0.001 mg/l | 0.001 mg/l |
| Toluene | <0.001 mg/l | 0.001 mg/l |
| Xylenes | <0.001 mg/l | 0.001 mg/l |
| Ethyl Benzene | <0.001 mg/l | 0.001 mg/l |
| Sb | 0.48 mg/l | 0.002 mg/l |
| As | <0.05 mg/l | 0.05 mg/l |
| Be | <0.01 mg/l | 0.01 mg/l |
| Cd | <0.01 mg/l | 0.01 mg/l |
| Cr | <0.05 mg/l | 0.05 mg/l |
| Cu | <0.03 mg/l | 0.03 mg/l |
| Pb | <0.05 mg/l | 0.05 mg/l |
| Hg | <0.002 mg/l | 0.002 mg/l |
| Ni | 0.08 mg/l | 0.06 mg/l |
| Se | 0.04 mg/l | 0.002 mg/l |
| Ag | <0.05 mg/l | 0.05 mg/l |
| Tl | <0.1 mg/l | 0.1 mg/l |
| Zn | 0.009 mg/l | 0.001 mg/l |

REFERENCES: "Test Methods for Evaluatin Solid Waste,-Physical/Chemical Methods", USEPA, SW 846, EMSL-Cincinnati, 1982.

An invoice for services is enclosed. Thank you for contacting Assaigai Laboratories.

Sincerely,

Jennifer V. Smith, Ph.D.
Laboratory Director



ASSAIGAI ANALYTICAL LABORATORIES



TO: Bloomfield Refinery
ATTN: Chris Hawley
PO Box 159
Bloomfield, NM 87413

DATE: 19 January 1987
2160

SAMPLE ID: MW-3

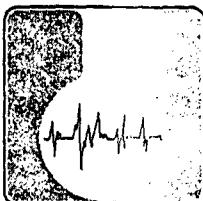
| ANALYTE | ANALYTICAL RESULTS | NOMINAL DETECTION LIMITS |
|---------------|--------------------|--------------------------|
| CN | 0.07 mg/l | 0.01 mg/l |
| Phenols | 0.012 mg/l | 0.001 mg/l |
| TOC | 12 mg/l | 0.1 mg/l |
| TDS | 4860 mg/l | 1 mg/l |
| Cl | 1290 mg/l | 1.0 mg/l |
| SO 4 | 2204 mg/l | 1.0 mg/l |
| Benzene | <0.001 mg/l | 0.001 mg/l |
| Toluene | <0.001 mg/l | 0.001 mg/l |
| Xylenes | <0.001 mg/l | 0.001 mg/l |
| Ethyl Benzene | <0.001 mg/l | 0.001 mg/l |
| Sb | 0.67 mg/l | 0.002 mg/l |
| As | <0.05 mg/l | 0.05 mg/l |
| Be | <0.01 mg/l | 0.01 mg/l |
| Cd | 0.11 mg/l | 0.01 mg/l |
| Cl | <0.05 mg/l | <0.05 mg/l |
| Cu | <0.03 mg/l | 0.03 mg/l |
| Pb | <0.05 mg/l | 0.05 mg/l |
| Hg | <0.002 mg/l | 0.002 mg/l |
| Ni | 0.10 mg/l | 0.06 mg/l |
| Se | 0.05 mg/l | 0.002 mg/l |
| Ag | <0.05 mg/l | 0.05 mg/l |
| Tl | <0.1 mg/l | 0.1 mg/l |
| Zn | 0.010 mg/l | 0.001 mg/l |

REFERENCES: "Test Methods for Evaluatin Solid Waste,-Physical/Chemical Methods", USEPA, SW 846, EMSL-Cincinnati, 1982.

An invoice for services is enclosed. Thank you for contacting Assaigai Laboratories.

Sincerely,

Jennifer V. Smith
Jennifer V. Smith, Ph.D.
Laboratory Director



ASSAIGAI ANALYTICAL LABORATORIES

TO: Bloomfield Refining
ATTN: Chris Hawley
PO Box 159
Bloomfield, NM 87413

DATE: 29 January 1987
2160 (revised)

SAMPLE ID: MW-4

ANALYTE

ANALYTICAL RESULTS

BASE NEUTRALS

| | |
|-----------------------------|-------------|
| Naphthalene | 0.036 mg/l |
| Acenaphthylene | <0.001 mg/l |
| Acenaphthene | 0.049 mg/l |
| Benzidine | <0.001 mg/l |
| 3,4-benzofluoranthene | <0.001 mg/l |
| Bis(2-chloroethoxy)methane | <0.001 mg/l |
| Bis(2-chloroethyl)ether | <0.001 mg/l |
| Bis(2-chloroisopropyl)ether | <0.001 mg/l |
| Bis(2-ethylhexyl)phthalate | <0.001 mg/l |
| 4-bromophenyl phenyl ether | <0.001 mg/l |
| Butylbenzyl phthalate | <0.001 mg/l |
| 2-chloronaphthalene | <0.001 mg/l |
| 4-chlorophenyl phenyl ether | <0.001 mg/l |
| 1,2-dichlorobenzene | <0.001 mg/l |
| 1,3-dichlorobenzene | <0.001 mg/l |
| 1,4-dichlorobenzene | <0.001 mg/l |
| 3,3-dichlorobenzidine | <0.001 mg/l |
| Diethyl phthalate | <0.001 mg/l |
| Dimethyl phthalate | <0.001 mg/l |
| Di-n-butyl phthalate | <0.001 mg/l |
| 2,4-dinitrotoluene | <0.001 mg/l |
| 2,6-dinitrotoluene | <0.001 mg/l |
| Di-n-octyl phthalate | <0.001 mg/l |
| 1,2-diphenylhydrazine | <0.001 mg/l |
| Fluoranthene | <0.001 mg/l |
| Fluorene | 0.023 mg/l |
| Hexachlorobenzene | <0.001 mg/l |
| Hexachlorobutadiene | <0.001 mg/l |
| Hexachlorocyclopentadiene | <0.001 mg/l |
| Hexachloroethane | <0.001 mg/l |
| Isophorone | <0.001 mg/l |
| Nitrobenzene | <0.001 mg/l |
| N-nitrosodimethylamine | <0.001 mg/l |
| N-nitrosodi-n-propylamine | <0.001 mg/l |
| N-nitrosodiphenylamine | <0.001 mg/l |
| 1,2,4-trichlorobenzene | <0.001 mg/l |

| | |
|------------------------|-------------|
| Phenanthrene | <0.001 mg/l |
| Anthracene | <0.001 mg/l |
| Pyrene | <0.001 mg/l |
| Benzo(a)anthracene | <0.001 mg/l |
| Chrysene | <0.001 mg/l |
| Benzo(k)fluoranthene | <0.001 mg/l |
| Benzo(a)pyrene | <0.001 mg/l |
| Dibenz(a,h)anthracene | <0.001 mg/l |
| Indeno(1,2,3-cd)pyrene | <0.001 mg/l |
| Benzo(g,h,i)perylene | <0.001 mg/l |

ACID PRIORITY POLLUTANTS

| | |
|------------------------|-------------|
| 2-chlorophenol | <0.001 mg/l |
| 2-nitrophenol | <0.001 mg/l |
| Phenol | <0.001 mg/l |
| 2,4-dimethyl phenol | <0.001 mg/l |
| 2,4-dichloro phenol | <0.001 mg/l |
| 2,4,6-trichloro phenol | <0.001 mg/l |
| P-chloro-m-cresol | <0.001 mg/l |
| 2,4 dinitro phenol | <0.001 mg/l |
| Pentachloro phenol | <0.001 mg/l |
| 4-nitro phenol | <0.001 mg/l |
| 4,6-dinitro-o-cresol | <0.001 mg/l |

VOLATILES

| | |
|--------------------------|-------------|
| Acrolein | <0.001 mg/l |
| Acrylonitrile | <0.001 mg/l |
| Chlorobenzene | <0.001 mg/l |
| Toluene | 1.78 mg/l |
| Benzene | 1.91 mg/l |
| Ethyl benzene | 4.48 mg/l |
| Tetrachloroethylene | <0.001 mg/l |
| Bromoform | <0.001 mg/l |
| Carbon tetrachloride | <0.001 mg/l |
| Chlorodibromomethane | <0.001 mg/l |
| Chloroethane | <0.001 mg/l |
| 2-chloroethylvinyl ether | <0.001 mg/l |
| Chloroform | <0.001 mg/l |
| Dichlorobromomethane | <0.001 mg/l |
| Dichlorodifluoromethane | <0.001 mg/l |
| 1,1-dichloroethane | <0.001 mg/l |
| 1,2-dichloroethane | <0.001 mg/l |
| 1,1-dichloroethylene | <0.001 mg/l |
| 1,2-dichloropropane | <0.001 mg/l |

| | |
|----------------------------|-------------|
| 1,2-dichloropropylene | <0.001 mg/l |
| Methyl bromide | <0.001 mg/l |
| Methyl chloride | <0.001 mg/l |
| Methylene chloride | <0.001 mg/l |
| 1,1,2,2-tetrachloroethane | <0.001 mg/l |
| 1,2-trans-dichloroethylene | <0.001 mg/l |
| 1,1,1-trichloroethane | <0.001 mg/l |
| 1,1,2-trichloroethane | <0.001 mg/l |
| Trichloroethylene | <0.001 mg/l |
| Vinyl chloride | <0.001 mg/l |

NOMINAL DETECTION LIMIT: 0.001 mg/l

| ANALYTE | ANALYTICAL RESULTS | NOMINAL DETECTION LIMITS |
|----------------------|--------------------|--------------------------|
| CN | <0.01 mg/l | 0.01 mg/l |
| Phenols | 0.096 mg/l | 0.001 mg/l |
| TOC | 170 mg/l | 0.1 mg/l |
| TDS | 2128 mg/l | 1 mg/l |
| Cl | 675 mg/l | 1.0 mg/l |
| SO ₄ | <1 mg/l | 1.0 mg/l |
| Sb | 0.40 mg/l | 0.01 mg/l |
| As | <0.05 mg/l | 0.05 mg/l |
| Be | <0.01 mg/l | 0.01 mg/l |
| Cd | <0.01 mg/l | 0.01 mg/l |
| Cr | <0.05 mg/l | 0.05 mg/l |
| Cu | <0.03 mg/l | 0.03 mg/l |
| Pb | <0.05 mg/l | 0.05 mg/l |
| Hg | <0.002 mg/l | 0.002 mg/l |
| Ni | <0.06 mg/l | 0.06 mg/l |
| Se | 0.03 mg/l | 0.002 mg/l |
| Ag | <0.05 mg/l | 0.05 mg/l |
| Tl | <0.1 mg/l | 0.1 mg/l |
| Zn | 0.040 mg/l | 0.001 mg/l |
| Ba | 2.3 mg/l | 0.05 mg/l |
| Fe | 18.6 mg/l | 0.3 mg/l |
| Mn | 5.70 mg/l | 0.05 mg/l |
| Al | 3.8 mg/l | 0.05 mg/l |
| B | 0.7 mg/l | 0.04 mg/l |
| Co | <0.05 mg/l | 0.05 mg/l |
| Mo | <0.01 mg/l | 0.01 mg/l |
| NO ₃ as N | <0.01 mg/l | 0.01 mg/l |
| F | 0.410 mg/l | 0.01 mg/l |

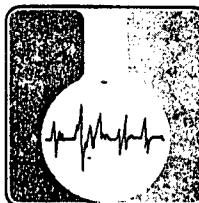
REFERENCE: "Test Methods for Evaluating Solid Waste, Physical/Chemical Methods," USEPA, SW 846, EMSL-Cincinnati, 1982.

An invoice for services is enclosed. Thank you for contacting Assaigai Laboratories.

Sincerely,



Jennifer V. Smith, Ph.D.
Laboratory Director



ASSAIGAI ANALYTICAL LABORATORIES



TO: Bloomfield Refinery
ATTN: Chris Hawley
PO Box 159
Bloomfield, NM 87413

DATE: 19 January 1987
2160

SAMPLE ID: MW-5

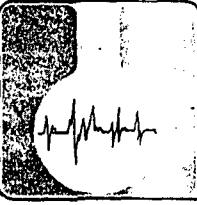
| ANALYTE | ANALYTICAL RESULTS | NOMINAL DETECTION LIMITS |
|----------------------|--------------------|--------------------------|
| CN | <0.01 mg/l | 0.01 mg/l |
| Phenols | 0.021 mg/l | 0.001 mg/l |
| TOC | 9 mg/l | 0.1 mg/l |
| TDS | 3788 mg/l | 1 mg/l |
| Cl | 1118 mg/l | 1.0 mg/l |
| SO ₄ | 1132 mg/l | 1.0 mg/l |
| Benzene | <0.001 mg/l | 0.001 mg/l |
| Toluene | <0.001 mg/l | 0.001 mg/l |
| Xylenes | <0.001 mg/l | 0.001 mg/l |
| Ethyl Benzene | <0.001 mg/l | 0.001 mg/l |
| Sb | 0.50 mg/l | 0.002 mg/l |
| As | <0.05 mg/l | 0.05 mg/l |
| Be | <0.01 mg/l | 0.01 mg/l |
| Cd | 0.01 mg/l | 0.01 mg/l |
| Cr | <0.05 mg/l | 0.05 mg/l |
| Cu | <0.03 mg/l | 0.03 mg/l |
| Pb | <0.05 mg/l | 0.05 mg/l |
| Hg | <0.002 mg/l | 0.002 mg/l |
| Ni | 0.07 mg/l | 0.06 mg/l |
| Se | 0.03 mg/l | 0.002 mg/l |
| Ag | <0.05 mg/l | 0.05 mg/l |
| Tl | <0.1 mg/l | 0.1 mg/l |
| Zn | 0.016 mg/l | 0.001 mg/l |
| Ba | 0.010 mg/l | 0.005 mg/l |
| Fe | <0.3 mg/l | 0.3 mg/l |
| Mn | <0.005 mg/l | 0.005 mg/l |
| Al | 4.34 mg/l | 0.05 mg/l |
| B | 0.24 mg/l | 0.004 mg/l |
| Co | <0.05 mg/l | 0.05 mg/l |
| Mo | 0.08 mg/l | 0.01 mg/l |
| NO ₃ as N | 36 mg/l | 0.1 mg/l |
| F | 0.580 mg/l | 0.01 mg/l |
| 1,2-dichloroethane | <0.001 mg/l | 0.001 mg/l |
| 1,1-dichloroethylene | <0.001 mg/l | 0.001 mg/l |
| Tetrachloroethylene | <0.001 mg/l | 0.001 mg/l |
| Trichloroethylene | <0.001 mg/l | 0.001 mg/l |
| CCl ₄ | <0.001 mg/l | 0.001 mg/l |

REFERENCES: "Test Methods for Evaluatin Solid Waste,-Physical/Chemical Methods", USEPA, SW 846, EMSL-Cincinnati, 1982.

An invoice for services is enclosed. Thank you for contacting Assaigai Laboratories.

Sincerely,

Jennifer V. Smith
Jennifer V. Smith, Ph.D.
Laboratory Director



ASSAIGAI ANALYTICAL LABORATORIES



TO: Bloomfield Refining
ATTN: Chris Hawley
PO Box 159
Bloomfield, NM 87413

DATE: 29 January 1987
2160 (revised)

SAMPLE ID: MW-7

ANALYTE

ANALYTICAL RESULTS

BASE NEUTRALS

| | |
|-----------------------------|-------------|
| Napthalene | <0.001 mg/l |
| Acenaphthylene | <0.001 mg/l |
| Acenaphthene | <0.001 mg/l |
| Benzidine | <0.001 mg/l |
| 3,4-benzofluoranthene | <0.001 mg/l |
| Bis(2-chloroethoxy)methane | <0.001 mg/l |
| Bis(2-chloroethyl)ether | <0.001 mg/l |
| Bis(2-chloroisopropyl)ether | <0.001 mg/l |
| Bis(2-ethylhexyl)phthalate | <0.001 mg/l |
| 4-bromophenyl phenyl ether | <0.001 mg/l |
| Butylbenzyl phthalate | <0.001 mg/l |
| 2-chloronaphthalene | <0.001 mg/l |
| 4-chlorophenyl phenyl ether | <0.001 mg/l |
| 1,2-dichlorobenzene | <0.001 mg/l |
| 1,3-dichlorobenzene | <0.001 mg/l |
| 1,4-dichlorobenzene | <0.001 mg/l |
| 3,3-dichlorobenzidine | <0.001 mg/l |
| Diethyl phthalate | <0.001 mg/l |
| Dimethyl phthalate | <0.001 mg/l |
| Di-n-butyl phthalate | <0.001 mg/l |
| 2,4-dinitrotoluene | <0.001 mg/l |
| 2,6-dinitrotoluene | <0.001 mg/l |
| Di-n-octyl phthalate | <0.001 mg/l |
| 1,2-diphenylhydrazine | <0.001 mg/l |
| Fluoranthene | <0.001 mg/l |
| Fluorene | <0.001 mg/l |
| Hexachlorobenzene | <0.001 mg/l |
| Hexachlorobutadiene | <0.001 mg/l |
| Hexachlorocyclopentadiene | <0.001 mg/l |
| Hexachloroethane | <0.001 mg/l |
| Isophorone | <0.001 mg/l |
| Nitrobenzene | <0.001 mg/l |
| N-nitrosodimethylamine | <0.001 mg/l |
| N-nitrosodi-n-propylamine | <0.001 mg/l |
| N-nitrosodiphenylamine | <0.001 mg/l |
| 1,2,4-trichlorobenzene | <0.001 mg/l |

| | |
|------------------------|-------------|
| Phenanthrene | <0.001 mg/l |
| Anthracene | <0.001 mg/l |
| Pyrene | <0.001 mg/l |
| Benzo(a)anthracene | <0.001 mg/l |
| Chrysene | 0.002 mg/l |
| Benzo(k)fluoranthene | 0.001 mg/l |
| Benzo(a)pyrene | <0.001 mg/l |
| Dibenz(a,h)anthracene | <0.001 mg/l |
| Indeno(1,2,3-cd)pyrene | <0.001 mg/l |
| Benzo(g,h,i)perylene | <0.001 mg/l |

ACID PRIORITY POLLUTANTS

| | |
|------------------------|-------------|
| 2-chlorophenol | <0.001 mg/l |
| 2-nitrophenol | <0.001 mg/l |
| Phenol | <0.001 mg/l |
| 2,4-dimethyl phenol | <0.001 mg/l |
| 2,4-dichloro phenol | <0.001 mg/l |
| 2,4,6-trichloro phenol | <0.001 mg/l |
| P-chloro-m-cresol | <0.001 mg/l |
| 2,4 dinitro phenol | <0.001 mg/l |
| Pentachloro phenol | <0.001 mg/l |
| 4-nitro phenol | <0.001 mg/l |
| 4,6-dinitro-o-cresol | <0.001 mg/l |

VOLATILES

| | |
|--------------------------|-------------|
| Acrolein | <0.001 mg/l |
| Acrylonitrile | <0.001 mg/l |
| Chlorobenzene | <0.001 mg/l |
| Toluene | <0.001 mg/l |
| Benzene | 0.009 mg/l |
| Ethyl benzene | <0.001 mg/l |
| Tetrachloroethylene | <0.001 mg/l |
| Bromoform | <0.001 mg/l |
| Carbon tetrachloride | <0.001 mg/l |
| Chlorodibromomethane | <0.001 mg/l |
| Chloroethane | <0.001 mg/l |
| 2-chloroethylvinyl ether | <0.001 mg/l |
| Chloroform | <0.001 mg/l |
| Dichlorobromomethane | <0.001 mg/l |
| Dichlorodifluoromethane | <0.001 mg/l |
| 1,1-dichloroethane | <0.001 mg/l |
| 1,2-dichloroethane | <0.001 mg/l |
| 1,1-dichloroethylene | <0.001 mg/l |
| 1,2-dichloropropane | <0.001 mg/l |

| | |
|----------------------------|-------------|
| 1,2-dichloropropylene | <0.001 mg/l |
| Methyl bromide | <0.001 mg/l |
| Methyl chloride | <0.001 mg/l |
| Methylene chloride | <0.001 mg/l |
| 1,1,2,2-tetrachloroethane | <0.001 mg/l |
| 1,2-trans-dichloroethylene | <0.001 mg/l |
| 1,1,1-trichloroethane | <0.001 mg/l |
| 1,1,2-trichloroethane | <0.001 mg/l |
| Trichloroethylene | <0.001 mg/l |
| Vinyl chloride | <0.001 mg/l |

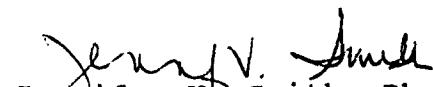
NOMINAL DETECTION LIMIT: 0.001 mg/l

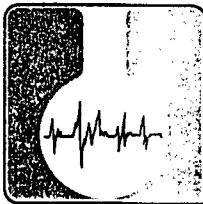
| ANALYTE | ANALYTICAL RESULTS | NOMINAL DETECTION LIMITS |
|-----------------|--------------------|--------------------------|
| CN | <0.01 mg/l | 0.01 mg/l |
| Phenols | 0.025 mg/l | 0.001 mg/l |
| TOC | 2 mg/l | 0.1 mg/l |
| TDS | 6940 mg/l | 1 mg/l |
| Cl | 29 mg/l | 1.0 mg/l |
| SO ₄ | 3630 mg/l | 1.0 mg/l |
| Sb | 0.83 mg/l | 0.01 mg/l |
| As | <0.05 mg/l | 0.05 mg/l |
| Be | <0.01 mg/l | 0.01 mg/l |
| Cd | 0.02 mg/l | 0.01 mg/l |
| Cr | 0.08 mg/l | 0.05 mg/l |
| Cu | <0.03 mg/l | 0.03 mg/l |
| Pb | 0.26 mg/l | 0.05 mg/l |
| Hg | <0.002 mg/l | 0.002 mg/l |
| Ni | 0.07 mg/l | 0.06 mg/l |
| Se | 0.09 mg/l | 0.002 mg/l |
| Ag | <0.05 mg/l | 0.05 mg/l |
| Tl | <0.1 mg/l | 0.1 mg/l |
| Zn | 0.017 mg/l | 0.001 mg/l |
| Xylenes | <0.001 mg/l | 0.001 mg/l |

REFERENCE: "Test Methods for Evaluating Solid Waste, Physical/Chemical Methods," USEPA, SW 846, EMSL-Cincinnati, 1982.

An invoice for services is enclosed. Thank you for contacting Assaigai Laboratories.

Sincerely,


Jennifer V. Smith, Ph.D.
Laboratory Director



ASSAIGAI ANALYTICAL LABORATORIES



TO: Bloomfield Refining
ATTN: Chris Hawley
PO Box 159
Bloomfield, NM 87413

DATE: 29 January 1987
2160 (revised)

SAMPLE ID: MW-8

ANALYTE

ANALYTICAL RESULTS

BASE NEUTRALS

| | |
|-----------------------------|-------------|
| Naphthalene | <0.001 mg/l |
| Acenaphthylene | <0.001 mg/l |
| Acenaphthene | <0.001 mg/l |
| Benzidine | <0.001 mg/l |
| 3,4-benzofluoranthene | <0.001 mg/l |
| Bis(2-chloroethoxy)methane | <0.001 mg/l |
| Bis(2-chloroethyl)ether | <0.001 mg/l |
| Bis(2-chloroisopropyl)ether | <0.001 mg/l |
| Bis(2-ethylhexyl)phthalate | <0.001 mg/l |
| 4-bromophenyl phenyl ether | <0.001 mg/l |
| Butylbenzyl phthalate | <0.001 mg/l |
| 2-chloronaphthalene | <0.001 mg/l |
| 4-chlorophenyl phenyl ether | <0.001 mg/l |
| 1,2-dichlorobenzene | <0.001 mg/l |
| 1,3-dichlorobenzene | <0.001 mg/l |
| 1,4-dichlorobenzene | <0.001 mg/l |
| 3,3-dichlorobenzidine | <0.001 mg/l |
| Diethyl phthalate | <0.001 mg/l |
| Dimethyl phthalate | <0.001 mg/l |
| Di-n-butyl phthalate | <0.001 mg/l |
| 2,4-dinitrotoluene | <0.001 mg/l |
| 2,6-dinitrotoluene | <0.001 mg/l |
| Di-n-octyl phthalate | <0.001 mg/l |
| 1,2-diphenylhydrazine | <0.001 mg/l |
| Fluoranthene | <0.001 mg/l |
| Fluorene | <0.001 mg/l |
| Hexachlorobenzene | <0.001 mg/l |
| Hexachlorobutadiene | <0.001 mg/l |
| Hexachlorocyclopentadiene | <0.001 mg/l |
| Hexachloroethane | <0.001 mg/l |
| Isophorone | <0.001 mg/l |
| Nitrobenzene | <0.001 mg/l |
| N-nitrosodimethylamine | <0.001 mg/l |
| N-nitrosodi-n-propylamine | <0.001 mg/l |
| N-nitrosodiphenylamine | <0.001 mg/l |
| 1,2,4-trichlorobenzene | <0.001 mg/l |

| | |
|------------------------|-------------|
| Phenanthrene | <0.001 mg/l |
| Anthracene | <0.001 mg/l |
| Pyrene | <0.001 mg/l |
| Benzo(a)anthracene | <0.001 mg/l |
| Chrysene | <0.001 mg/l |
| Benzo(k)fluoranthene | <0.001 mg/l |
| Benzo(a)pyrene | <0.001 mg/l |
| Dibenz(a,h)anthracene | <0.001 mg/l |
| Indeno(1,2,3-cd)pyrene | <0.001 mg/l |
| Benzo(g,h,i)perylene | <0.001 mg/l |

ACID PRIORITY POLLUTANTS

| | |
|------------------------|-------------|
| 2-chlorophenol | <0.001 mg/l |
| 2-nitrophenol | <0.001 mg/l |
| Phenol | <0.001 mg/l |
| 2,4-dimethyl phenol | <0.001 mg/l |
| 2,4-dichloro phenol | <0.001 mg/l |
| 2,4,6-trichloro phenol | <0.001 mg/l |
| P-chloro-m-cresol | <0.001 mg/l |
| 2,4 dinitro phenol | <0.001 mg/l |
| Pentachloro phenol | <0.001 mg/l |
| 4-nitro phenol | <0.001 mg/l |
| 4,6-dinitro-o-cresol | <0.001 mg/l |

VOLATILES

| | |
|--------------------------|-------------|
| Acrolein | <0.001 mg/l |
| Acrylonitrile | <0.001 mg/l |
| Chlorobenzene | <0.001 mg/l |
| Toluene | <0.001 mg/l |
| Benzene | <0.001 mg/l |
| Ethyl benzene | <0.001 mg/l |
| Tetrachloroethylene | <0.001 mg/l |
| Bromoform | <0.001 mg/l |
| Carbon tetrachloride | <0.001 mg/l |
| Chlorodibromomethane | <0.001 mg/l |
| Chloroethane | <0.001 mg/l |
| 2-chloroethylvinyl ether | <0.001 mg/l |
| Chloroform | <0.001 mg/l |
| Dichlorobromomethane | <0.001 mg/l |
| Dichlorodifluoromethane | <0.001 mg/l |
| 1,1-dichloroethane | <0.001 mg/l |
| 1,2-dichloroethane | <0.001 mg/l |
| 1,1-dichloroethylene | <0.001 mg/l |
| 1,2-dichloropropane | <0.001 mg/l |

| | |
|----------------------------|-------------|
| 1,2-dichloropropylene | <0.001 mg/l |
| Methyl bromide | <0.001 mg/l |
| Methyl chloride | <0.001 mg/l |
| Methylene chloride | <0.001 mg/l |
| 1,1,2,2-tetrachloroethane | <0.001 mg/l |
| 1,2-trans-dichloroethylene | <0.001 mg/l |
| 1,1,1-trichloroethane | <0.001 mg/l |
| 1,1,2-trichloroethane | <0.001 mg/l |
| Trichloroethylene | <0.001 mg/l |
| Vinyl chloride | <0.001 mg/l |

NOMINAL DETECTION LIMIT: 0.001 mg/l

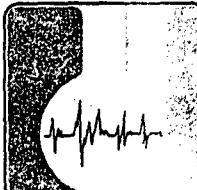
| ANALYTE | ANALYTICAL RESULTS | NOMINAL DETECTION LIMITS |
|-----------------|--------------------|--------------------------|
| CN | 0.10 mg/l | 0.01 mg/l |
| Phenols | 0.042 mg/l | 0.001 mg/l |
| TOC | 8 mg/l | 0.1 mg/l |
| TDS | 3450 mg/l | 1 mg/l |
| Cl | 913 mg/l | 1.0 mg/l |
| SO ₄ | 1270 mg/l | 1.0 mg/l |
| Sb | 0.67 mg/l | 0.01 mg/l |
| As | <0.05 mg/l | 0.05 mg/l |
| Be | <0.01 mg/l | 0.01 mg/l |
| Cd | <0.01 mg/l | 0.01 mg/l |
| Cr | <0.05 mg/l | 0.05 mg/l |
| Cu | <0.03 mg/l | 0.03 mg/l |
| Pb | <0.05 mg/l | 0.05 mg/l |
| Hg | <0.002 mg/l | 0.002 mg/l |
| Ni | 0.43 mg/l | 0.06 mg/l |
| Se | 0.04 mg/l | 0.002 mg/l |
| Ag | <0.05 mg/l | 0.05 mg/l |
| Tl | <0.1 mg/l | 0.1 mg/l |
| Zn | 0.016 mg/l | 0.001 mg/l |
| Xylenes | <0.001 mg/l | 0.001 mg/l |

REFERENCE: "Test Methods for Evaluating Solid Waste, Physical/Chemical Methods," USEPA, SW 846, EMSL-Cincinnati, 1982.

An invoice for services is enclosed. Thank you for contacting Assaigai Laboratories.

Sincerely,

Jennifer V. Smith, Ph.D.
Laboratory Director



ASSAIGAI
ANALYTICAL
LABORATORIES

TO: Bloomfield Refining
ATTN: Chris Hawley
PO Box 159
Bloomfield, NM 87413

DATE: 29 January 1987
2160 (revised)

SAMPLE ID: MW-9

ANALYTE

ANALYTICAL RESULTS

BASE NEUTRALS

| | |
|-----------------------------|-------------|
| Naphthalene | 0.029 mg/l |
| Acenaphthylene | <0.001 mg/l |
| Acenaphthene | <0.001 mg/l |
| Benzidine | <0.001 mg/l |
| 3,4-benzofluoranthene | <0.001 mg/l |
| Bis(2-chloroethoxy)methane | <0.001 mg/l |
| Bis(2-chloroethyl)ether | <0.001 mg/l |
| Bis(2-chloroisopropyl)ether | <0.001 mg/l |
| Bis(2-ethylhexyl)phthalate | <0.001 mg/l |
| 4-bromophenyl phenyl ether | <0.001 mg/l |
| Butylbenzyl phthalate | <0.001 mg/l |
| 2-chloronaphthalene | <0.001 mg/l |
| 4-chlorophenyl phenyl ether | <0.001 mg/l |
| 1,2-dichlorobenzene | <0.001 mg/l |
| 1,3-dichlorobenzene | <0.001 mg/l |
| 1,4-dichlorobenzene | <0.001 mg/l |
| 3,3-dichlorobenzidine | <0.001 mg/l |
| Diethyl phthalate | <0.001 mg/l |
| Dimethyl phthalate | <0.001 mg/l |
| Di-n-butyl phthalate | <0.001 mg/l |
| 2,4-dinitrotoluene | <0.001 mg/l |
| 2,6-dinitrotoluene | <0.001 mg/l |
| Di-n-octyl phthalate | <0.001 mg/l |
| 1,2-diphenylhydrazine | <0.001 mg/l |
| Fluoranthene | <0.001 mg/l |
| Fluorene | <0.001 mg/l |
| Hexachlorobenzene | <0.001 mg/l |
| Hexachlorobutadiene | <0.001 mg/l |
| Hexachlorocyclopentadiene | <0.001 mg/l |
| Hexachloroethane | <0.001 mg/l |
| Isophorone | <0.001 mg/l |
| Nitrobenzene | <0.001 mg/l |
| N-nitrosodimethylamine | <0.001 mg/l |
| N-nitrosodi-n-propylamine | <0.001 mg/l |
| N-nitrosodiphenylamine | <0.001 mg/l |
| 1,2,4-trichlorobenzene | <0.001 mg/l |

| | |
|------------------------|-------------|
| Phenanthrene | <0.001 mg/l |
| Anthracene | <0.001 mg/l |
| Pyrene | <0.001 mg/l |
| Benzo(a)anthracene | <0.001 mg/l |
| Chrysene | <0.001 mg/l |
| Benzo(k)fluoranthene | <0.001 mg/l |
| Benzo(a)pyrene | <0.001 mg/l |
| Dibenz(a,h)anthracene | <0.001 mg/l |
| Indeno(1,2,3-cd)pyrene | <0.001 mg/l |
| Benzo(g,h,i)perylene | <0.001 mg/l |

ACID PRIORITY POLLUTANTS

| | |
|------------------------|-------------|
| 2-chlorophenol | <0.001 mg/l |
| 2-nitrophenol | <0.001 mg/l |
| Phenol | 0.133 mg/l |
| 2,4-dimethyl phenol | <0.001 mg/l |
| 2,4-dichloro phenol | <0.001 mg/l |
| 2,4,6-trichloro phenol | <0.001 mg/l |
| P-chloro-m-cresol | <0.001 mg/l |
| 2,4 dinitro phenol | <0.001 mg/l |
| Pentachloro phenol | <0.001 mg/l |
| 4-nitro phenol | <0.001 mg/l |
| 4,6-dinitro-o-cresol | <0.001 mg/l |

VOLATILES

| | |
|--------------------------|-------------|
| Acrolein | <0.001 mg/l |
| Acrylonitrile | <0.001 mg/l |
| Chlorobenzene | <0.001 mg/l |
| Toluene | 0.754 mg/l |
| Benzene | 1.49 mg/l |
| Ethyl benzene | 0.504 mg/l |
| Tetrachloroethylene | <0.001 mg/l |
| Bromoform | <0.001 mg/l |
| Carbon tetrachloride | <0.001 mg/l |
| Chlorodibromomethane | <0.001 mg/l |
| Chloroethane | <0.001 mg/l |
| 2-chloroethylvinyl ether | <0.001 mg/l |
| Chloroform | <0.001 mg/l |
| Dichlorobromomethane | <0.001 mg/l |
| Dichlorodifluoromethane | <0.001 mg/l |
| 1,1-dichloroethane | <0.001 mg/l |
| 1,2-dichloroethane | <0.001 mg/l |
| 1,1-dichloroethylene | <0.001 mg/l |
| 1,2-dichloropropane | <0.001 mg/l |

| | |
|----------------------------|-------------|
| 1,2-dichloropropylene | <0.001 mg/l |
| Methyl bromide | <0.001 mg/l |
| Methyl chloride | <0.001 mg/l |
| Methylene chloride | <0.001 mg/l |
| 1,1,2,2-tetrachloroethane | <0.001 mg/l |
| 1,2-trans-dichloroethylene | <0.001 mg/l |
| 1,1,1-trichloroethane | <0.001 mg/l |
| 1,1,2-trichloroethane | <0.001 mg/l |
| Trichloroethylene | <0.001 mg/l |
| Vinyl chloride | <0.001 mg/l |

NOMINAL DETECTION LIMIT: 0.001 mg/l

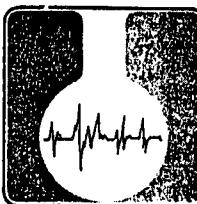
| ANALYTE | ANALYTICAL RESULTS | NOMINAL DETECTION LIMITS |
|-----------------|--------------------|--------------------------|
| CN | <0.01 mg/l | 0.01 mg/l |
| Phenols | 0.160 mg/l | 0.001 mg/l |
| TOC | 275 mg/l | 0.1 mg/l |
| TDS | 1684 mg/l | 1 mg/l |
| Cl | 109 mg/l | 1.0 mg/l |
| SO ₄ | 20 mg/l | 1.0 mg/l |
| Sb | 0.40 mg/l | 0.01 mg/l |
| As | <0.05 mg/l | 0.05 mg/l |
| Be | <0.01 mg/l | 0.01 mg/l |
| Cd | <0.01 mg/l | 0.01 mg/l |
| Cr | <0.05 mg/l | 0.05 mg/l |
| Cu | <0.03 mg/l | 0.03 mg/l |
| Pb | <0.05 mg/l | 0.05 mg/l |
| Hg | <0.002 mg/l | 0.002 mg/l |
| Ni | 0.16 mg/l | 0.06 mg/l |
| Se | 0.03 mg/l | 0.002 mg/l |
| Ag | <0.05 mg/l | 0.05 mg/l |
| Tl | <0.1 mg/l | 0.1 mg/l |
| Zn | 0.011 mg/l | 0.001 mg/l |
| Xylenes | <0.001 mg/l | 0.001 mg/l |

REFERENCE: "Test Methods for Evaluating Solid Waste, Physical/Chemical Methods," USEPA, SW 846, EMSL-Cincinnati, 1982.

An invoice for services is enclosed. Thank you for contacting Assaigai Laboratories.

Sincerely,

Jennifer V. Smith
Jennifer V. Smith, Ph.D.
Laboratory Director



ASSAIGAI ANALYTICAL LABORATORIES



TO: Bloomfield Refining
ATTN: Chris Hawley
PO Box 159
Bloomfield, NM 87413

DATE: 19 January 1987
2160

SAMPLE ID: MW-10

ANALYTE

ANALYTICAL RESULTS

BASE NEUTRALS

| | |
|-----------------------------|-------------|
| Naphthalene | 0.004 mg/l |
| Acenaphthylene | <0.001 mg/l |
| Acenaphthene | <0.001 mg/l |
| Benzidine | <0.001 mg/l |
| 3,4-benzofluoranthene | <0.001 mg/l |
| Bis(2-chloroethoxy)methane | <0.001 mg/l |
| Bis(2-chloroethyl)ether | <0.001 mg/l |
| Bis(2-chloroisopropyl)ether | <0.001 mg/l |
| Bis(2-ethylhexyl)phthalate | <0.001 mg/l |
| 4-bromophenyl phenyl ether | <0.001 mg/l |
| Butylbenzyl phthalate | <0.001 mg/l |
| 2-chloronaphthalene | <0.001 mg/l |
| 4-chlorophenyl phenyl ether | <0.001 mg/l |
| 1,2-dichlorobenzene | <0.001 mg/l |
| 1,3-dichlorobenzene | <0.001 mg/l |
| 1,4-dichlorobenzene | <0.001 mg/l |
| 3,3-dichlorobenzidine | <0.001 mg/l |
| Diethyl phthalate | <0.001 mg/l |
| Dimethyl phthalate | <0.001 mg/l |
| Di-n-butyl phthalate | <0.001 mg/l |
| 2,4-dinitrotoluene | <0.001 mg/l |
| 2,6-dinitrotoluene | <0.001 mg/l |
| Di-n-octyl phthalate | <0.001 mg/l |
| 1,2-diphenylhydrazine | <0.001 mg/l |
| Fluoranthene | <0.001 mg/l |
| Fluorene | <0.001 mg/l |
| Hexachlorobenzene | <0.001 mg/l |
| Hexachlorobutadiene | <0.001 mg/l |
| Hexachlorocyclopentadiene | <0.001 mg/l |
| Hexachloroethane | <0.001 mg/l |
| Isophorone | <0.001 mg/l |
| Nitrobenzene | <0.001 mg/l |
| N-nitrosodimethylamine | <0.001 mg/l |
| N-nitrosodi-n-propylamine | <0.001 mg/l |
| N-nitrosodiphenylamine | <0.001 mg/l |
| 1,2,4-trichlorobenzene | <0.001 mg/l |

| | |
|------------------------|-------------|
| Phenanthrene | <0.001 mg/l |
| Anthracene | <0.001 mg/l |
| Pyrene | <0.001 mg/l |
| Benzo(a)anthracene | <0.001 mg/l |
| Chrysene | <0.001 mg/l |
| Benzo(k)fluoranthene | <0.001 mg/l |
| Benzo(a)pyrene | <0.001 mg/l |
| Dibenz(a,h)anthracene | <0.001 mg/l |
| Indeno(1,2,3-cd)pyrene | <0.001 mg/l |
| Benzo(g,h,i)perylene | <0.001 mg/l |

ACID PRIORITY POLLUTANTS

| | |
|------------------------|-------------|
| 2-chlorophenol | <0.001 mg/l |
| 2-nitrophenol | <0.001 mg/l |
| Phenol | <0.001 mg/l |
| 2,4-dimethyl phenol | <0.001 mg/l |
| 2,4-dichloro phenol | <0.001 mg/l |
| 2,4,6-trichloro phenol | <0.001 mg/l |
| P-chloro-m-cresol | <0.001 mg/l |
| 2,4 dinitro phenol | <0.001 mg/l |
| Pentachloro phenol | <0.001 mg/l |
| 4-nitro phenol | <0.001 mg/l |
| 4,6-dinitro-o-cresol | <0.001 mg/l |

VOLATILES

| | |
|--------------------------|-------------|
| Acrolein | <0.001 mg/l |
| Acrylonitrile | <0.001 mg/l |
| Chlorobenzene | <0.001 mg/l |
| Toluene | 7.4 mg/l |
| Benzene | 14.4 mg/l |
| Ethyl benzene | 0.030 mg/l |
| Tetrachloroethylene | <0.001 mg/l |
| Bromoform | <0.001 mg/l |
| Carbon tetrachloride | <0.001 mg/l |
| Chlorodibromomethane | <0.001 mg/l |
| Chloroethane | <0.001 mg/l |
| 2-chloroethylvinyl ether | <0.001 mg/l |
| Chloroform | <0.001 mg/l |
| Dichlorobromomethane | <0.001 mg/l |
| Dichlorodifluoromethane | <0.001 mg/l |
| 1,1-dichloroethane | <0.001 mg/l |
| 1,2-dichloroethane | <0.001 mg/l |
| 1,1-dichloroethylene | <0.001 mg/l |
| 1,2-dichloropropane | <0.001 mg/l |

| | |
|----------------------------|-------------|
| 1,2-dichloropropylene | <0.001 mg/l |
| Methyl bromide | <0.001 mg/l |
| Methyl chloride | <0.001 mg/l |
| Methylene chloride | <0.001 mg/l |
| 1,1,2,2-tetrachloroethane | <0.001 mg/l |
| 1,2-trans-dichloroethylene | <0.001 mg/l |
| 1,1,1-trichloroethane | <0.001 mg/l |
| 1,1,2-trichloroethane | <0.001 mg/l |
| Trichloroethylene | <0.001 mg/l |
| Vinyl chloride | <0.001 mg/l |

NOMINAL DETECTION LIMIT: 0.001 mg/l

| ANALYTE | ANALYTICAL RESULTS | NOMINAL DETECTION LIMITS |
|-----------------|--------------------|--------------------------|
| CN | <0.01 mg/l | 0.01 mg/l |
| Phenols | 0.055 mg/l | 0.001 mg/l |
| TOC | 114 mg/l | 0.1 mg/l |
| TDS | 3272 mg/l | 1 mg/l |
| Cl | 457 mg/l | 1.0 mg/l |
| SO ₄ | 10 mg/l | 1.0 mg/l |
| Sb | 0.56 mg/l | 0.002 mg/l |
| As | <0.05 mg/l | 0.05 mg/l |
| Be | 0.04 mg/l | 0.01 mg/l |
| Cd | <0.01 mg/l | 0.01 mg/l |
| Cr | <0.05 mg/l | 0.05 mg/l |
| Cu | <0.03 mg/l | 0.03 mg/l |
| Pb | <0.05 mg/l | 0.05 mg/l |
| Hg | <0.002 mg/l | 0.002 mg/l |
| Ni | <0.06 mg/l | 0.06 mg/l |
| Se | 0.03 mg/l | 0.002 mg/l |
| Ag | <0.05 mg/l | 0.05 mg/l |
| Tl | <0.1 mg/l | 0.1 mg/l |
| Zn | 0.010 mg/l | 0.001 mg/l |
| Xylenes | <0.001 mg/l | 0.001 mg/l |

REFERENCE: "Test Methods for Evaluating Solid Waste, Physical/Chemical Methods," USEPA, SW 846, EMSL-Cincinnati, 1982.

An invoice for services is enclosed. Thank you for contacting Assaigai Laboratories.

Sincerely,

Jennifer V. Smith
Jennifer V. Smith, Ph.D.
Laboratory Director

TO: Bloomfield Refinery

0695

Page 3 of 3

NOMINAL DETECTION LIMITS

| | |
|---------|------------|
| Phenols | 0.002 mg/l |
| Benzene | 0.001 mg/l |
| Toluene | 0.001 mg/l |

Detection limits for Volatiles, and Base/Neutrals are
all 0.001 mg/l

An invoice for services is enclosed. Thank you for contacting Assaigai
Laboratories.

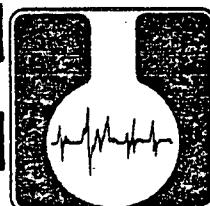
Sincerely,



Jennifer V. Smith, Ph.D.
Laboratory Director

APPENDIX B

**SURFACE WATER SAMPLING RESULTS
HAMMOND DITCH**



ASSAIGAI ANALYTICAL LABORATORIES



TO: Bloomfield Refinery
Attn: Chris Hawley
P.O. Box 159
Bloomfield NM 87413

HAMMOND DITCH

DATE: 22 May 1986
0660 completed
Page 1 of 4

SAMPLE DATE: 4/22/86

ANALYTE

| | SAMPLE ID/ ANALYTICAL RESULTS | |
|--------------------------|-------------------------------|--------------------------|
| | NEAR SULLIVAN ROAD HSRD 5 | NEAR API PONDS HAPI 5 |
| CN | <0.01 mg/l | <0.01 mg/l |
| Phenols | 0.002 mg/l | 0.002 mg/l |
| Sb | <0.2 mg/l | <0.2 mg/l |
| As | <0.050 mg/l | <0.050 mg/l |
| Be | <0.01 mg/l | <0.01 mg/l |
| Cd | <0.010 mg/l | <0.010 mg/l |
| Cr | <0.050 mg/l | <0.050 mg/l |
| Cu | <0.03 mg/l | <0.03 mg/l |
| Pb | <0.050 mg/l | <0.050 mg/l |
| Hg | <0.002 mg/l | <0.002 mg/l |
| Ni | <0.06 mg/l | <0.06 mg/l |
| Se | <0.010 mg/l | <0.010 mg/l |
| Ag | <0.050 mg/l | <0.050 mg/l |
| Tl | <0.01 mg/l | <0.01 mg/l |
| Zn | <0.01 mg/l | <0.01 mg/l |
| Acrolein | ND | ND |
| Acrylonitrile | ND | ND |
| Benzene | 0.006 mg/l | ND |
| Bromoform | ND | ND |
| Carbon Tetrachloride | ND | ND |
| Chlorobenzene | ND | ND |
| Chlorodibromomethane | ND | ND |
| Chloroethane | ND | ND |
| 2-Chloroethylvinyl Ether | ND | ND |
| Chloroform | ND | ND |
| Dichlorogromomethane | ND | ND |
| 1,1-Dichloroethane | ND | ND |
| 1,2-Dichloroethane | ND | ND |
| 1,1-Dichloroethylene | ND | ND |
| 1,2-Dichloropropane | ND | ND |
| 1,2-Dichloropropylene | ND | ND |
| Ethylbenzene | ND | ND |

TO: Bloomfield Refinery

0660

Page 2 of 4

ANALYTE

SAMPLE ID/ ANALYTICAL RESULTS

HSRD 5

HAPI 5

| | | |
|-----------------------------|------------|----|
| Methyl Bromide | ND | ND |
| Methyl Chloride | ND | ND |
| Methylene Chloride | ND | ND |
| 1,1,2,2-Tetrachloroethane | ND | ND |
| Tetrachloroethylene | ND | ND |
| Toluene | 0.003 mg/l | ND |
| 1,2-Transdichloroethylene | ND | ND |
| 1,1,1-Trichloroethane | ND | ND |
| 1,1,2-Trichloroethane | ND | ND |
| Trichloroethylene | ND | ND |
| Vinyl Chloride | ND | ND |
| Acid Compounds | | |
| 2-chlorophenol | ND | ND |
| 2,4-dichlorophenol | ND | ND |
| 2,4-dimethylphenol | ND | ND |
| 4,6-dinitro-o-cresol | ND | ND |
| 2,4-dinitrophenol | ND | ND |
| 2-nitrophenol | ND | ND |
| 4-nitrophenol | ND | ND |
| p-chloro-m-cresol | ND | ND |
| pentachlorophenol | ND | ND |
| Phenol | ND | ND |
| 2,4,6-trichlorophenol | ND | ND |
| Base Neutrals | | |
| Acenaphthene | ND | ND |
| Acenaphthylene | ND | ND |
| Anthracene | 0.006 mg/l | ND |
| Benzidine | ND | ND |
| Benzo(a)anthracene | 0.003 mg/l | ND |
| Benzo(a)pyrene | ND | ND |
| 3,4-benzofluoranthene | ND | ND |
| Benzo(ghi)perylene | ND | ND |
| Benzo(k)fluoranthene | ND | ND |
| Bis(2-chloroethoxy)methane | ND | ND |
| Bis(2-chloroethyl)ether | ND | ND |
| Bis(2-chloroisopropyl)ether | ND | ND |
| Bis(2-ethylhexyl)phthalate | ND | ND |
| 4-bromophenyl phenyl ether | ND | ND |
| Butylbenzyl phthalate | ND | ND |
| 2-chloronaphthalene | ND | ND |
| 4-chlorophenyl phenyl ether | ND | ND |
| Chrysene | 0.005 mg/l | ND |

TO: Bloomfield Refinery

0660 completed
Page 3 of 4

ANALYTE

SAMPLE ID/ ANALYTICAL RESULTS

HSRD 5

HAPI 5

| | | |
|---------------------------|------------|------------|
| Dibenzo(a,h)anthracene | ND | ND |
| 1,2-Dichlorobenzene | ND | ND |
| 1,3-Dichlorobenzene | ND | ND |
| 1,4-Dichlorobenzene | ND | ND |
| 3,3-Dichlorobenzidine | ND | ND |
| Diethyl phthalate | ND | ND |
| Dimethyl phthalate | ND | ND |
| Din-n-butyl phthalate | ND | ND |
| 2,4-dinitrotoluene | ND | ND |
| 2,6-dinitrotoluene | ND | ND |
| Di-n-octyl phthalate | ND | ND |
| 1,2-diphenylhydrazine | ND | ND |
| Fluoranthene | ND | 0.001 mg/l |
| Fluorene | ND | ND |
| Hexachlorobenzene | ND | ND |
| Hexachlorobutadiene | ND | ND |
| Hexachlorocyclopentadiene | ND | ND |
| Hexachloroethane | ND | ND |
| Indeno(1,2,3-cd)pyrene | ND | ND |
| Isophorone | ND | ND |
| Naphthalene | 0.013 mg/l | ND |
| Nitrobenzene | ND | ND |
| N-nitrosodimethylamine | ND | ND |
| N-nitrosodi-n-propylamine | ND | ND |
| N-nitrosodiphenylamine | ND | ND |
| Phenanthrene | 0.007 mg/l | ND |
| Pyrene | 0.008 mg/l | ND |
| 1,2,4-trichlorobenzene | ND | ND |

ND = None Detected

REFERENCE: "Test Methods for Evaluating Solid Waste Chemical/Physical Methods", USEPA, SW 846, EMSL-Cincinnati, 1982.

TO: Bloomfield Refinery

0660 completed
Page 4 of 4

NOMINAL DETECTION LIMITS

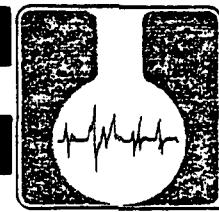
| | |
|--------------|------------|
| CN | 0.01 mg/l |
| Phenols | 0.002 mg/l |
| Benzene | 0.001 mg/l |
| Toluene | 0.001 mg/l |
| Xylenes | 0.001 mg/l |
| Ethylbenzene | 0.001 mg/l |
| Sb | 0.2 mg/l |
| As | 0.050 mg/l |
| Be | 0.010 mg/l |
| Cd | 0.010 mg/l |
| Cr | 0.050 mg/l |
| Cu | 0.03 mg/l |
| Pb | 0.050 mg/l |
| Hg | 0.002 mg/l |
| Ni | 0.06 mg/l |
| Se | 0.010 mg/l |
| Ag | 0.050 mg/l |
| Tl | 0.01 mg/l |
| Zn | 0.01 mg/l |

Detection limits for Volatiles, Acid Compounds, and Base/Neutrals are
all 0.001 mg/l

An invoice for services is enclosed. Thank you for contacting Assaigai
Laboratories.

Sincerely,


Jennifer V. Smith, Ph.D.
Laboratory Director



ASSAIGAI ANALYTICAL LABORATORIES

TO: Bloomfield Refinery
Attn: Chris Hawley
P.O. Box 159
Bloomfield NM 87413

DATE: 22 May 1986

0695

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SAMPLE DATE: 4/28/86

HAMMOND DITCH

ANALYTE

SAMPLE ID/ ANALYTICAL RESULTS

U4

NEAR API WASTE PONDS

0.003 mg/l

D4

NEAR SULLIVAN ROAD

0.002 mg/l

U6A

D6A

| | | |
|----------------------------------|----|----|
| Acrolein | ND | ND |
| Acrylonitrile | ND | ND |
| Benzene | ND | ND |
| <u>Bromoform</u> | ND | ND |
| Carbon Tetrachloride | ND | ND |
| Chlorobenzene | ND | ND |
| Chlorodibromomethane | ND | ND |
| Chloroethane | ND | ND |
| 2-Chloroethylvinyl Ether | ND | ND |
| Chloroform | ND | ND |
| Dichlorogromomethane | ND | ND |
| 1,1-Dichloroethane | ND | ND |
| 1,2-Dichloroethane | ND | ND |
| 1,1-Dichloroethylene | ND | ND |
| 1,2-Dichloropropane | ND | ND |
| 1,2-Dichloropropylene | ND | ND |
| Ethylbenzene | ND | ND |
| Methyl Bromide | ND | ND |
| Methyl Chloride | ND | ND |
| Methylene Chloride | ND | ND |
| 1,1,2,2-Tetrachloroethane | ND | ND |
| Tetrachloroethylene | ND | ND |
| Toluene | ND | ND |
| <u>1,2-Transdichloroethylene</u> | ND | ND |
| 1,1,1-Trichloroethane | ND | ND |
| 1,1,2-Trichloroethane | ND | ND |
| Trichloroethylene | ND | ND |
| Vinly Chloride | ND | ND |
| Base Neutrals | | |
| Acenaphthene | ND | ND |
| Acenphthylene | ND | ND |
| Anthracene | ND | ND |
| <u>Benzidine</u> | ND | ND |

TO: Bloomfield Refinery

0695

Page 2 of 3

ANALYTE

SAMPLE ID/ ANALYTICAL RESULTS

U6A

D6A

| | | |
|-----------------------------|----|----|
| Benzo(a)anthracene | ND | ND |
| Benzo(a)pyrene | ND | ND |
| 3,4-benzofluoranthene | ND | ND |
| Benzo(ghi)perylene | ND | ND |
| Benzo(k)fluoranthene | ND | ND |
| Bis(2-chloroethoxy)methane | ND | ND |
| Bis(2-chloroethyl)ether | ND | ND |
| Bis(2-chloroisopropyl)ether | ND | ND |
| Bis(2-ethylhexyl)phthalate | ND | ND |
| 4-bromophenyl phenyl ether | ND | ND |
| Butylbenzyl phthalate | ND | ND |
| 2-chloronaphthalene | ND | ND |
| 4-chlorophenyl phenyl ether | ND | ND |
| Chrysene | ND | ND |
| <hr/> | | |
| Dibenzo(a,h)anthracene | ND | ND |
| 1,2-Dichlorobenzene | ND | ND |
| 1,3-Dichlorobenzene | ND | ND |
| 1,4-Dichlorobenzene | ND | ND |
| 3,3-Dichlorobenzidine | ND | ND |
| Diethyl phthalate | ND | ND |
| Dimethyl phthalate | ND | ND |
| Din-n-butyl phthalate | ND | ND |
| 2,4-dinitrotoluene | ND | ND |
| 2,6-dinitrotoluene | ND | ND |
| Di-n-octyl phthalate | ND | ND |
| 1,2-diphenylhydrazine | ND | ND |
| Fluoranthene | ND | ND |
| <hr/> | | |
| Fluorene | ND | ND |
| Hexachlorobenzene | ND | ND |
| Hexachlorobutadiene | ND | ND |
| Hexachlorocyclopentadiene | ND | ND |
| Hexachloroethane | ND | ND |
| Indeno(1,2,3-cd)pyrene | ND | ND |
| Isophorone | ND | ND |
| <u>Naphthalene</u> | ND | ND |
| <hr/> | | |
| Nitrobenzene | ND | ND |
| N-nitrosodimethylamine | ND | ND |
| N-nitrosodie-n-propylamine | ND | ND |
| N-nitrosodiphenylamine | ND | ND |
| <u>Phenanthrene</u> | ND | ND |
| <hr/> | | |
| Pyrene | ND | ND |
| <hr/> | | |
| 1,2,4-trichlorobenzene | ND | ND |

ND = None Detected

REFERENCE: "Test Methods for Evaluating Solid Waste Chemical/Physical Methods", USEPA, SW 846, EMSL-Cincinnati, 1982

APPENDIX C

**EPA AMBIENT STANDARDS AND CRITERIA
FOR SUPERFUND REMEDIAL SITES**

APPENDIX C

**EPA AMBIENT STANDARDS AND CRITERIA
FOR SUPERFUND REMEDIAL SITES**

| Chemical | Applicable or Relevant Requirements | | Other Criteria, Advisories, and Guidance | |
|----------------------------|---|---|--|---|
| | Clean Air Act, MCLs (mg/L unless otherwise noted) | RAQS (ug/m ³) | Clean Water Act, Water Quality Criteria for Human Health— Fish and Drinking Water | Clean Water Act, Water Quality Criteria for Human Health— Drinking Water Only ^a |
| Acenaphthene | | | 20 ug/L (organoleptic) ^b | 20 ug/L (organoleptic) |
| Acrolein | | | 320 ug/L | 560 ug/L |
| Acrylonitrile | | | 0 (58 ng/L) ^c | 0 (63 ng/L) |
| Aldrin | | | 0 (0.076 ng/L) | 0 (1.2 ng/L) |
| Antimony | 0.05 | | 146 ug/L | 146 ug/L |
| Arsenic | | | 0 (2.2 ng/L) | 0 (2.5 ng/L) |
| Asbestos | 1.0 | | 0 (30,000 fibers/L) | 0 (30,000 fibers/L) |
| Barium | | | 0 (0.66 ug/L) | 0 (0.67 ug/L) |
| Benzene | | | 0 (0.12 ng/L) | 0 (0.15 ng/L) |
| Benzidine | | | 0 (3.7 ng/L) | 0 (3.9 ng/L) |
| Beryllium | | | 10 ug/L | 10 ug/L |
| Cadmium | 0.01 | 40,000 (1-hour) ^d 10,000 (8-hour) | 0 (0.4 ug/L) | 0 (0.42 ug/L) |
| Carbon monoxide | | | 0 (0.46 ng/L) | 0 (22 ng/L) |
| Carbon tetrachloride | | | 0 (0.72 ng/L) | 0 (21 ng/L) |
| Chlordane | | | 38 ug/L | 160 ug/L |
| Chlorinated benzenes | | | 74 ug/L | 570 ug/L |
| Hexachlorobenzene | | | Insufficient data | Insufficient data |
| 1,2,4,5-Tetrachlorobenzene | | | 488 ug/L | 488 ug/L |
| Pentachlorobenzene | | | 0 (0.94 ug/L) | 0 (0.94 ug/L) |
| Trichlorobenzene | | | 18.4 mg/L | 19 mg/L |
| Monochlorobenzene | | | 0 (0.6 ug/L) | 0 (0.6 ug/L) |
| 1,1,1-Trichloroethane | | | 0 (0.17 ug/L) | 0 (0.17 ug/L) |
| 1,1,2-Trichloroethane | | | 0 (1.9 ug/L) | 0 (2.4 ug/L) |
| 1,1,2,2-Tetrachloroethane | | | Insufficient data | Insufficient data |
| Heptachloroethane | | | Insufficient data | Insufficient data |
| 1,1-Dichloroethane | | | Insufficient data | Insufficient data |
| 1,1,1,2-Tetrachloroethane | | | Insufficient data | Insufficient data |
| Pentachloroethane | | | Insufficient data | (continued) |

| Chemical | Applicable or Relevant Requirements | | Other Criteria, Advisories, and Guidance | | | |
|--|---|--|--|--|--|--|
| | Safe Drinking Water Act, MCLs ($\mu\text{g/L}$, unless otherwise noted) | Clean Air Act, NAAQS ($\mu\text{g/m}^3$) | Clean Water Act, Water Quality Criteria for Human Health—Fish and Drinking Water | Clean Water Act, Water Quality Criteria for Human Health—Adjusted for Drinking Water Only* | Safe Drinking Water Act, Health Advisories ($\mu\text{g/L}$) | Chronic 1-day 10-day Chronic (longer term) |
| Chlorinated naphthalenes | | | Inadequate data | Inadequate data | | |
| Chlorinated phenols | | | | | | |
| 3-Monochlorophenol | | | 0.1 $\mu\text{g/L}$ (organoleptic) | 0.1 $\mu\text{g/L}$ (organoleptic) | 0.1 $\mu\text{g/L}$ (organoleptic) | |
| 4-Monochlorophenol | | | 0.1 $\mu\text{g/L}$ (organoleptic) | 0.1 $\mu\text{g/L}$ (organoleptic) | 0.1 $\mu\text{g/L}$ (organoleptic) | |
| 2,3-Dichlorophenol | | | 0.06 $\mu\text{g/L}$ (organoleptic) | 0.04 $\mu\text{g/L}$ (organoleptic) | 0.04 $\mu\text{g/L}$ (organoleptic) | |
| 2,3,5-Dichlorophenol | | | 0.5 $\mu\text{g/L}$ (organoleptic) | 0.5 $\mu\text{g/L}$ (organoleptic) | 0.5 $\mu\text{g/L}$ (organoleptic) | |
| 2,6-Dichlorophenol | | | 0.2 $\mu\text{g/L}$ (organoleptic) | 0.2 $\mu\text{g/L}$ (organoleptic) | 0.2 $\mu\text{g/L}$ (organoleptic) | |
| 3,4-Dichlorophenol | | | 0.1 $\mu\text{g/L}$ (organoleptic) | 0.3 $\mu\text{g/L}$ (organoleptic) | 0.3 $\mu\text{g/L}$ (organoleptic) | |
| 2,3,4,6-Tetrachlorophenol | | | 1.0 $\mu\text{g/L}$ (organoleptic) | 1.0 $\mu\text{g/L}$ (organoleptic) | 1.0 $\mu\text{g/L}$ (organoleptic) | |
| 2,4,5-Trichlorophenol | | | 2600 $\mu\text{g/L}$ | 2600 $\mu\text{g/L}$ | 0 (1.8 $\mu\text{g/L}$) | |
| 2,4,6-Trichloropheno ¹ | | | 0 (1.2 $\mu\text{g/L}$) | 1800 $\mu\text{g/L}$ (organoleptic) | 1800 $\mu\text{g/L}$ (organoleptic) | |
| 2-Methyl-4-chlorophenol | | | 3000 $\mu\text{g/L}$ (organoleptic) | 3000 $\mu\text{g/L}$ (organoleptic) | 3000 $\mu\text{g/L}$ (organoleptic) | |
| 3-Methyl-4-chloropheno ¹ | | | 20 $\mu\text{g/L}$ (organoleptic) | 20 $\mu\text{g/L}$ (organoleptic) | 20 $\mu\text{g/L}$ (organoleptic) | |
| Chlorophenoxy ² | | | | | | |
| 2,4-Dichlorophenoxyacetic acid (2,4-D) | 0.1 | | | | | |
| 2,4,5-Trichlorophenoxy-propionic acid (2,4,5-TP) | 0.01 | | | | | |
| Chloroalyl ethers | | | | | | |
| bis-(Chloromethyl) ether | | | 0 (0.0038 $\mu\text{g/L}$) | 0 (0.0039 $\mu\text{g/L}$) | 0 (0.0039 $\mu\text{g/L}$) | |
| bis-(2-Chloroethyl) ether | | | 0 (30 $\mu\text{g/L}$) | 0 (30 $\mu\text{g/L}$) | 0 (30 $\mu\text{g/L}$) | |
| bis-(2-Chloroisopropyl) ether | | | 34.7 $\mu\text{g/L}$ | 34.7 $\mu\text{g/L}$ | 34.7 $\mu\text{g/L}$ | |
| Chloroform | 0.1e | | 0 (0.19 $\mu\text{g/L}$) | 0 (0.19 $\mu\text{g/L}$) | 0 (0.19 $\mu\text{g/L}$) | |
| 2-Chlorophenol | 0.05 | | 0.1 $\mu\text{g/L}$ (organoleptic) | 0.1 $\mu\text{g/L}$ (organoleptic) | 0.1 $\mu\text{g/L}$ (organoleptic) | |
| Chromium Cr+6 Cr+3 | | | 50 $\mu\text{g/L}$ | 50 $\mu\text{g/L}$ | 50 $\mu\text{g/L}$ | |
| Copper | | | 170 $\mu\text{g/L}$ | 170 $\mu\text{g/L}$ | 179 $\mu\text{g/L}$ | |
| Cyanide | | | 1 $\mu\text{g/L}$ (organoleptic) | 1 $\mu\text{g/L}$ (organoleptic) | 1 $\mu\text{g/L}$ (organoleptic) | |
| DNT | | | 200 $\mu\text{g/L}$ | 200 $\mu\text{g/L}$ | 200 $\mu\text{g/L}$ | |
| Dichlorobenzenes (all isomers) | | | 400 $\mu\text{g/L}$ | 400 $\mu\text{g/L}$ | 470 $\mu\text{g/L}$ | |
| Dichlorobenzenes | | | 0 (10.3 $\mu\text{g/L}$) | 0 (10.3 $\mu\text{g/L}$) | 0 (20.7 $\mu\text{g/L}$) | |
| 1,1-Dichloroethylene | | | 0 (33 $\mu\text{g/L}$) | 0 (33 $\mu\text{g/L}$) | 0 (33 $\mu\text{g/L}$) | |
| 1,2-Dichloroethylene | | | Inadequate data | Inadequate data | 0.27 | (continued) |
| | | | | | 0.6 (cis isomer) 0.07 (trans isomer) | |

| Chemical | Applicable or Relevant Requirements | | Other Criteria, Advisories, and Guidance | | |
|--------------------------------------|--|---|--|---|---|
| | Clean Air Act ^t , MCLs (ug/L, unless otherwise noted) | Clean Air Act ^t , RAQS (ug/m ³) | Clean Water Act, Water Quality Criteria for Human Health— Fish and Drinking Water | Clean Water Act, Water Quality Criteria for Human Health— Adjusted for Drinking Water Only ^a | Safe Drinking Water Act, Health Advisories (ug/L) |
| Dichloromethane | | | See Halomethanes 3.09 ug/L | See Halomethanes 3.09 ug/L | 1.3 1.3 0.15 |
| 2,4-Dichlorophenol | | | | | Inadequate data 87 ug/L |
| Dichloropropane/ Dichloropropenes | | | | | 0 (0.071 ug/L) |
| Dichloropropanes | | | | | 400 ug/L (organoleptic) |
| Dieldrin | | | | | 0 (0.11 ug/L) |
| 2,4-Diethylphenol | | | | | 5.68 0.568 |
| 2-n-Dinitrotoluene | | | | | 0 (42 ug/L) |
| p-Dioxane | | | | | 74 ug/L |
| 1,2-Diphenylhydrazine | | | | | 138 ug/L |
| Endosulfan | 0.0002 | | | | 1 ug/L |
| Endrin | | | | | 1.4 ug/L |
| Ethylbenzene | | | | | 2.4 ug/L |
| Ethyleneglycol | | | | | 19.0 |
| Formaldehyde | | | | | 0.03 |
| Fluoranthene | 1.4-2.4 | | 42 ug/L | 188 ug/L | |
| Fluoride | | | | | Inadequate data 0 (0.19 ug/L) |
| Halothanes | | | | | 0 (0.28 ug/L) |
| Halomethanes | | | | | 0 (0.45 ug/L) |
| Heptachlorobutadiene | | | | | |
| Heptachloroclohexanes | | | | | |
| Lindane (99% gamma-HCH) | 0.004 | | | | |
| alpha-HCH | | | 0 (9.2 ug/L) | 0 (13 ug/L) | |
| beta-HCH | | | 0 (16.3 ug/L) | 0 (23.2 ug/L) | |
| gamma-HCH | | | 0 (18.6 ug/L) | 0 (26.4 ug/L) | |
| delta-HCH | | | | | Inadequate data |
| epsilon-HCH | | | | | Inadequate data |
| Technical-HCH | | | | | Inadequate data |
| Hexachlorocyclopentadiene | | | 0 (12.3 ug/L) | 0 (17.4 ug/L) | |
| n-Hexane | | | 206 ug/L | 206 ug/L | |
| Hydrocarbons (non-methane) | | | | | |
| Isophorone | | | | | |
| Kerosene/fuel oil no. 2 | | | | | |
| | | | 5.2 ug/L | 5.2 ug/L | 0.358 |
| | | | | | 0.23 |

(continued)

| Applicable or Relevant Requirements | | Other Criteria, Advisories, and Guidance | | | |
|---|---|---|---|--|---|
| Chemical | Clean Air Act, MCLs (mg/L unless otherwise noted) | Clean Air Act, MAQs (ug/m ³) | Clean Water Act, Water Quality Criteria for Human Health-- Fish and Drinking Water | Clean Water Act, Water Quality Criteria for Human Health-- Adjusted for Drinking Water Only* | Safe Drinking Water Act, Health Advisories (mg/L) |
| | | | 1-day | 10-day | Chronic (longer term) |
| Lead | 0.05 | 1.5 (90-day) | 50 ug/L | 50 ug/L | |
| Mercury | 0.002 | | 144 ng/L | 10 ug/L | |
| Methoxychlor | 0.1 | | | | Inadequate data 15.4 ug/L |
| Methyl Ethyl Ketone | | | | | Inadequate data 13.4 ug/L |
| Naphthalene | | | | | Inadequate data 19.8 ug/L |
| Nickel | 10.0 | 100 (1-year) ^h | | | |
| Nitrate (as N) | | | | | |
| Nitrobenzene | | | | | |
| Nitrogen dioxide | | | | | |
| Nitrophenoole | | | | | |
| 2,4-Dinitro- <i>o</i> -cresol | | | | | |
| Dinitrophenol | | | | | |
| Mononitrophenol | | | | | |
| Trinitrophenol | | | | | |
| Nitrosamines | | | | | |
| n-Nitrosodimethylamine | | | 0 (1.4 ng/L) | 0 (1.4 ng/L) | |
| n-Nitrosodiethylamine | | | 0 (0.8 ng/L) | 0 (0.8 ng/L) | |
| <i>o</i> -Nitrosodi- <i>n</i> -butylamine | | | 0 (6.4 ng/L) | 0 (6.4 ng/L) | |
| <i>n</i> -Nitrosodiphenylamine | | | 0 (4.9 ug/L) | 0 (7.0 ug/L) | |
| <i>n</i> -Nitrosopyrrolidine | | | 0 (16 ng/L) | 0 (16 ng/L) | |
| Ozone | | | | | |
| Particulate Matter | | | | | |
| Pentachlorophenol | | | 235 (1-hour) ^d | 1.01 mg/L | |
| Phenol | | | 260 (24-hour) ^d | 3.5 ug/L | |
| Phthalate esters | | | 75 (24-hour) ⁱ | | 3.5 ug/L |
| Dimethylphthalate | | | | 313 ug/L | |
| Diethylphthalate | | | | 350 ug/L | |
| Dibutylphthalate | | | | 434 ug/L | |
| Di-2-ethylhexyl-phthalate | | | | 44 ug/L | |
| Polychlorinated biphenyls (PCBs) | | | | 15 ug/L | |
| Polynuclear aromatic hydrocarbons (PAHs) | | | | 21 ug/L | |
| | | | | | |

(continued)

| Applicable or Relevant Requirements | | Other Criteria, Advisories, and Guidance | | | |
|--------------------------------------|---|--|--|---|---|
| Chemical | Water Act, MCLs (ug/L unless otherwise noted) | Clean Air Act, NAAQS (ug/m ³) | Clean Water Act, Water Quality Criteria for Human Health— Fish and Drinking Water | Clean Water Act, Water Quality Criteria for Human Health— Adjusted for Drinking Water Only ^a | Safe Drinking Water Act, Health Advisories (ug/L) 1-day 10-day Chronic (longer term) |
| Radionuclides | | | | | |
| Radium-226 and 228 | 5 pCi/L | | | 0 (0.000013 ug/L) | 0 (0.00016 ug/L) |
| Gross alpha activity | 15 pCi/L | | | 0 (0.8 ug/L) | 0 (0.88 ug/L) |
| Tritium | 20,000 pCi/L | | | 17.8 ug/L | 2.3 |
| Strontium-90 | 0 pCi/L | | | 15 ug/L | 2.2 |
| Other man-made | | | | 0 (25.8 ug/L) | 0.34 |
| Selenium | 0.01 | | 10 ug/L | 0 (0.71 ug/L) | 0 |
| Silver | 0.05 | | 50 ug/L | 0 (2.7 ug/L) | 0.075 |
| Sulfur dioxide | | 365 (24-hour) ^b 80 (1-year) ^b | | 0 (2.0 ug/L) | 21.5 |
| 2,3,7,8-TCDD | | | | (2.0 ug/L) | 2.0 |
| Tetrachloroethylene | | | | 0 (0.00016 ug/L) | 0.2 |
| Thallium | | | | 0 (0.8 ug/L) | 0.175 |
| Toluene | | | | 14.3 ug/L | 0.62 |
| Toxaphene | 0.005 | | | 0 (0.71 ug/L) | 0.075 |
| Trichloroethylene | | | | 0 (2.7 ug/L) | 0.075 |
| Trihalomethanes (total) ^c | 0.1 | | | 0 (2.0 ug/L) | 0.075 |
| Vinyl chloride | | | | 0 (0.8 ug/L) | 0.075 |
| Xylenes | | | | 0 (0.8 ug/L) | 0.075 |
| Zinc | | | | 0 (0.8 ug/L) | 0.075 |
| | | | | 5 ug/L (organoleptic) | 5 ug/L (organoleptic) |

^aThese adjusted criteria, for drinking water ingestion only, were derived from published EPA Water Quality Criteria (45 FR 79318-79379, November 28, 1980) for combined fish and drinking water ingestion and for fish ingestion alone. These adjusted values are not official EPA Water Quality Criteria, but may be appropriate for Superfund sites with contaminated ground water. In the derivation of these values, intake was assumed to be 2 liters/day for drinking water and 6.5 grams/day for fish; human body weight was assumed to be 70 kilograms.

^bCriteria designated as organoleptic are based on taste and odor effects, not human health effects. Health-based Water Quality Criteria are not available for these chemicals.

^cThe criterion for all carcinogens is zero; the concentration given in parentheses corresponds to a carcinogenic risk of 10^{-6} . Water Quality Criteria documents present concentrations resulting in risks from 10^{-6} to 10^{-5} . To obtain concentrations corresponding to risks of 10^{-6} and 10^{-5} , the 10^{-6} concentrations should be multiplied by 100 and 10, respectively. To obtain concentrations corresponding to risk of 10^{-7} , 10^{-6} concentrations should be divided by 10.

(continued)

^dAnnual maximum concentration not to be exceeded more than once per year.

^eChloroform is one of four trihalomethanes whose sum concentration must be less than 0.1 mg/L.

^fAs a guide in devising implementation plans for achieving oxidant standards.

^gSeven-day health advisory for benzene and benzo(a)pyrene in kerosene, respectively.

^hAnnual arithmetic mean concentration.

ⁱAnnual geometric mean concentration.

^jActivity corresponding to total body or any internal organ dose of 4 mrem/year.

^kTotal trihalomethanes refers to the sum concentration of chloroform, bromodichloromethane, dibromochloromethane, and bromoform.