United States Army
Fort Monmouth, New Jersey

Underground Storage Tank Closure and Site Investigation Report

Building 290
Main Post

NJDEP UST Registration No. 081533-64
NJDEP Closure Approval No. C-93-3179

May 2000

VOLUME 1 OF 3

UNDERGROUND STORAGE TANK
CLOSURE AND SITE INVESTIGATION REPORT

BUILDING 290

MAIN POST
NJDEP UST REGISTRATION NO. 081533-64
NJDEP CLOSURE APPROVAL NO. C-93-3179

MAY 2000

PROJECT NO.: 09-5004-08
CONTRACT NO.: DACA51-94-D-0014

PREPARED FOR:

UNITED STATES ARMY, FORT MONMOUTH, NEW JERSEY
DIRECTORATE OF PUBLIC WORKS
BUILDING 167
FORT MONMOUTH, NJ 07703

PREPARED BY:

ATC ASSOCIATES INC.
BROMLEY CORPORATE CENTER
THREE TERRI LANE
BURLINGTON, NEW JERSEY 08016

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EXECUTIVE SUMMARY

UST Closure

On September 12, 1994, a fiberglass underground storage tank (UST) was closed by removal in accordance with the New Jersey Department of Environmental Protection (NJDEP) Closure Approval No. C-93-3179 at U.S. Army Fort Monmouth, Fort Monmouth, New Jersey. The UST, NJDEP Registration No. 081533-64, was located immediately adjacent to Building 290 in the Main Post area of U.S. Army, Fort Monmouth. UST No. 081533-64 was a 2,000-gallon No. 2 diesel oil UST. The UST fill port was located directly above the tank. The tank closure was performed by Cleaning Up The Environment Inc. (CUTE Inc.).

Site Assessment - Soil

The site assessment was performed by U.S. Army personnel in accordance with the NJDEP *Technical Requirements for Site Remediation* (N.J.A.C. 7:26E) and the NJDEP *Field Sampling Procedures Manual*. Soils surrounding the tank were screened visually and with air monitoring equipment for evidence of contamination. Following removal, the UST was inspected for corrosion holes. No holes were noted in the UST, however, evidence of potentially contaminated soils was observed surrounding the tank.

On September 12, 1994, following the removal of the UST, approximately 40 cubic yards of potentially contaminated soil was removed from the excavation due to visual contamination.

On September 13, 1994, following removal of approximately 10 cubic yards of potentially contaminated soil, post-excavation soil samples A, B, C, D, E, F, and DUP A were collected from a total of a total of six (6) locations along the sidewalls of the excavation. The samples were collected at a depth of 5.5 feet below ground surface (bgs). Sample H was collected along the former piping length of the excavation, which was approximately 15 feet in length. The piping samples were collected at a depth of 5.5 feet bgs. All samples were analyzed for total petroleum hydrocarbons (TPHC).

On September 21, 1994, due to elevated TPHC levels, a sample was collected from the north sidewall of the excavation in the vicinity of sample A, and was analyzed for volatile organic compounds plus 15 tentatively identified compounds (VOCs). The sample was collected at a depth of 5.5 feet bgs.

Findings - Soil

Post-excavation soil samples (samples C, D, and H) collected from the UST excavation and from below piping associated with the former UST at Building 290 contained TPHC concentrations below the NJDEP residential direct contact total organic contaminants soil cleanup criteria of

10,000 milligrams per kilogram (mg/kg) (N.J.A.C. 7:26D and revisions dated February 3, 1994). Samples C, D, and H, contained levels of TPHC ranging in concentration from 57.3 mg/kg to 730.0 mg/kg. Samples E, and F contained TPHC concentrations of 3,110.0 mg/kg and 9,670.0 mg/kg, respectively. Samples A, B, and DUP A contained levels of TPHC ranging in concentration from 10,400.0 mg/kg to 16,200.0 mg/kg, which exceeded the NJDEP soil cleanup criteria for 10,000 mg/kg. Post-excavation soil sample A, collected on September 21, 1994, contained methylene chloride at 0.46 mg/kg, and 2-butanone at 1.5 mg/kg. No other compounds were detected.

Based on field screening of subsurface soils, the DPW has concluded that an historical discharge was associated with the UST and associated piping. On September 13, 1994, a spill was reported to the NJDEP "Hotline" for UST No. 081533-64 and was assigned Spill Case No. 94-9-13-1503-57.

Site Restoration

Following receipt of all post-excavation soil sampling results, the excavation was backfilled to grade with a combination of uncontaminated excavated soil and certified clean fill. The excavation site was then restored to its original condition.

Site Assessment - Groundwater

One shallow overburden monitoring well (MW-1) was installed at the Building 290 area on July 15, 1994. It was installed approximately 30 feet south off the southeastern corner of Building 290 in the assumed downgradient direction of the former excavation. It was screened in the 2.0- to 12.5- foot depth interval, across the water table, which is approximately 3 feet below grade surface.

A second shallow overburden monitoring well (MW-2) was installed at the Building 290 area on August 16, 1995. It was installed approximately 8 feet north off the northwestern corner of Building 290 in the assumed downgradient direction of the former excavation. It was screened in the 1.5- to 11.5- foot depth interval, across the water table, which is approximately 4 feet below grade surface.

Monitoring well MW-1 was sampled on November 8, 1994, November 29, 1994, and December 18, 1995, and analyzed for volatile organic compounds with xylenens (VOCs), tertiary butyl alcohol (TBA), methyl teriary butyl ether (MTBE), and total lead.

Monitoring well MW-2 was sampled on November 18, 1995, and analyzed for volatile organic compounds with xylenes (VOCs), tertiary butyl alcohol, methyl tertiary butyl ether (MTBE), and total lead.

All sampling and analyses were performed in accordance with the NJDEP *Field Sampling Procedures Manual*, and the *Technical Requirements for Site Remediation*, N.J.A.C. 7:26E (*Technical Requirements*).

Findings - Groundwater

The sample collected from MW-1 on November 8, 1994, contained lead at 17.0 micrograms per liter (ug/l). This exceeded the Ground Water Quality Criteria (GWQC) for lead of 10 ug/l. All other groundwater analytical results were either below the detection limit or in compliance with the New Jersey Groundwater Quality Criteria (GWQC).

The sample collected on November 29, 1994, contained lead at 3.0 ug/l, which complies with the GWQC for lead. No other compounds were detected.

The sample collected on December 18, 1995, contained methylene chloride at 1.1 ug/l, and cis-1,2-Dichloroethene at 1.1 ug/l. The sample also contained aluminum at 2,200.0 ug/l, barium at 55.0 ug/l, calcium at 4,100.0 ug/l, copper at 60.0 ug/l, iron at 6,400.0 ug/l, potassium at 5,000.0 ug/l, magnesium at 4,500.0 ug/l, manganese at 23.0 ug/l, sodium at 9,200.0 ug/l, and zinc at 180.0 ug/l. This exceeds the GWQC for alumium of 200.0 ug/l, and iron of 300.0 ug/l. The trip blank and the field blank contained methylene chloride at 1.4 ug/l. No other compounds were detected.

The sample collected from MW-2 on November 28, 1995 contained methylene chloride at 2.3 ug/l, cis-1,2-Dichloroethene at 1.3 ug/l, and chloroform at 0.50 ug/l. The sample also contained silver at 440.0 ug/l, aluminum at 360.0 ug/l, barium at 74.0 ug/l, calcium at 5,300.0 ug/l, iron at 1,100.0 ug/l, magnesium at 4,700.0 ug/l, manganese at 30.0 ug/l, sodium at 11,000.0 ug/l, lead at 1.5 ug/l, and zinc at 100.0 ug/l. This exceeds the GWQC for aluminum of 200.0 ug/l, and iron of 300.0 ug/l. The trip blank and the field blank contained methylene chloride at 0.70 ug/l, and 0.60 ug/l, respectively. The field blank also contained manganese at 32.0 ug/l, lead at 2.8 ug/l, and zinc at 27.0 ug/l. No other compounds were detected.

No product or sheen was observed in MW-1 or MW-2 on any of the sampling dates. The depth to the water table in MW-1 on November 8, 1994 was 8.14 feet below grade, 7.15 feet on November 29, 1994, and 3.06 feet below grade on December 18, 1995. The depth to the water table in MW-2 was 3.18 feet below grade on November 28, 1995.

Site Assessment Quality Assurance

The sampling and laboratory analysis conducted during the site assessment were performed in accordance with Section 7:26E-2.1 of the *Technical Requirements*.

Conclusions and Recommendations

Based on the post-excavation soil sampling results, soils with TPHC concentrations exceeding the NJDEP soil cleanup criteria for total organic contaminants of 10,000 mg/kg, do exist in the former location of the UST or associated piping.

The groundwater sample collected on November 8, 1994, contained a lead concentration which exceeded the New Jersey GWQC of 10 ug/l. However, lead was detected below the GWQC during the second round of groundwater sampling on November 29, 1994. Based on the analytical results of the groundwater samples collected on November 29, 1994, November 28, 1995, and December 18, 1995, groundwater quality at the Building 290 UST closure site complies with the New Jersey GWQC.

No further action is proposed in regard to the closure and site assessment of UST No. 081533-64 at Building 290.

1.0 UNDERGROUND STORAGE TANK DECOMMISSIONING ACTIVITIES

1.1 OVERVIEW

One underground storage tank (UST), New Jersey Department of Environmental Protection (NJDEP) Registration No. 081533-64, was closed at Building 290 at U.S. Army Fort Monmouth, Fort Monmouth, New Jersey on September 12, 1994. Refer to site location map on Figure 1. This report presents the results of the DPW's implementation of the UST Decommissioning/Closure Plan submitted to the NJDEP on June 28, 1993. The plan was approved on August 26, 1993 and assigned TMS No. C-93-3179. The UST was a fiberglass 2,000-gallon tank containing No. 2 diesel oil.

Decommissioning activities for UST No. 081533-64 complied with all applicable Federal, State and Local laws and ordinances in effect at the date of decommissioning. These laws included but were not limited to: N.J.A.C. 7:14B-1 et seq., N.J.A.C. 5:23-1 et seq., and Occupational Safety and Health Administration (OSHA) 1910.146 & 1910.120. All permits including but not limited to the NJDEP-approved Decommissioning/Closure Plan were posted onsite for inspection. CUTE Inc., the contractor that conducted the decommissioning activities, is registered and certified by the NJDEP for performing UST closure activities. Closure of UST No. 081533-64 proceeded under the approval of the NJDEP Bureau of Underground Storage Tanks (NJDEP-BUST). The NJDEP-BUST closure approval and signed certifications for UST No. 081533-64 are included in Appendices A and B, respectively.

Based on field screening of subsurface soils, the DPW has concluded that an historical discharge was associated with the UST and associated piping. On September 13, 1994, a spill was reported to the NJDEP "Hotline" for UST No. 081533-64 and was assigned Spill Case No. 94-9-13-1503-57.

This UST Closure and Site Investigation Report has been prepared by Smith Technology Corporation, to assist the United States Army Directorate of Public Works (DPW) in complying with the NJDEP Bureau of Underground Storage Tanks (NJDEP-BUST) regulations. The applicable NJDEP-BUST regulations at the date of closure were the *Interim Closure Requirements for Underground Storage Tank Systems* (N.J.A.C. 7:14B-1 et seq. September 1990 and revisions dated November 1, 1991).

This report was prepared using information required at the time of closure. Section 1 of this UST Closure and Site Investigation Report provides a summary of the UST decommissioning activities. Section 2 of this report describes the site investigation activities. Conclusions and recommendations, including the results of the soil sampling investigation, are presented in the final section of this report.

1.2 Site Description

Building 290 is located in the central portion of the Main Post area of Fort Monmouth, as shown on Figure 1. UST No. 081533-64 was located southwest of Building 290 and appurtenant piping ran approximately 15 feet northeast from the excavation to Building 290. The fill port area was located directly above the tank. A site map is provided on Figure 2.

1.2.1 Geological/Hydrogeological Setting

The following is a description of the geological/hydrogeological setting of the area surrounding Building 290. Included is a description of the regional geology of the area surrounding Fort Monmouth as well as descriptions of the local geology and hydrogeology of the Main Post area.

Regional Geology

Monmouth County lies within the New Jersey Section of the Atlantic Coastal Plain physiographic province. The Main Post, Charles Wood, and the Evans areas are located in what may be referred to as the Outer Coastal Plain subprovince, or the Outer Lowlands.

In general, New Jersey Coastal Plain formations consist of a seaward-dipping wedge of unconsolidated deposits of clay, silt, and gravel. These formations typically strike northeast-southwest with a dip ranging from 10 to 60 feet per mile and were deposited on Precambrian and lower Paleozoic rocks (Zapecza, 1989). These sediments, predominantly derived from deltaic, shallow marine, and continental shelf environments, date from Cretaceous through the Quaternary Periods. The mineralogy ranges from quartz to glauconite.

The formations record several major transgressive/regressive cycles and contain units which are generally thicker to the southeast and reflect a deeper water environment. Over 20 regional geologic units are present within the sediments of the Coastal Plain. Regressive, upward coarsening deposits are usually aquifers (e.g., Englishtown and Kirkwood Formations, and the Cohansey Sand) while the transgressive deposits act as confining units (e.g., the Merchantville, Marshalltown, and Navesink Formations). The individual thicknesses for these units vary greatly (i.e., from several feet to several hundred feet). The Coastal Plain deposits thicken to the southeast from the Fall Line to greater than 6,500 feet in Cape May County (Brown and Zapecza, 1990).

Local Geology

Based on the regional geologic map (Jablonski, 1968), the Cretaceous age Red Bank and Tinton Sands outcrop at the Main Post area. The Red Bank sand conformably overlies the Navesink Formation and dips to the southeast at 35 feet per mile. The upper member (Shrewsbury) of the Red Bank sand is a yellowish-gray to reddish brown clayey, medium-to-

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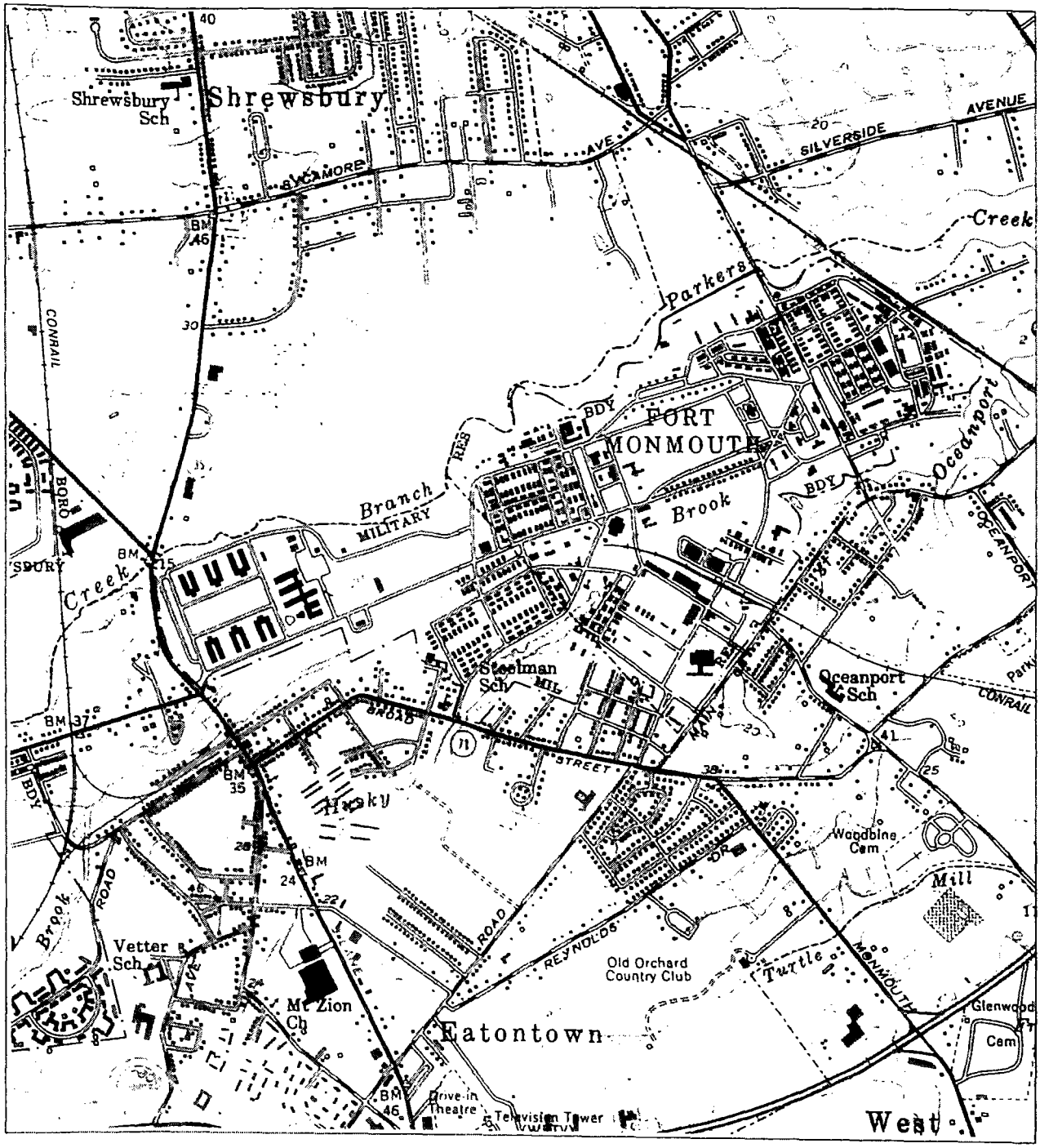
Fort Monmouth, New Jersey



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SCALE
0 2000 FT.

NEW JERSEY

QUADRANGLE LOCATION

Source: Long Branch, New Jersey Quadrangle

Project No. 09-5004-14

Figure 1
Site Location Map

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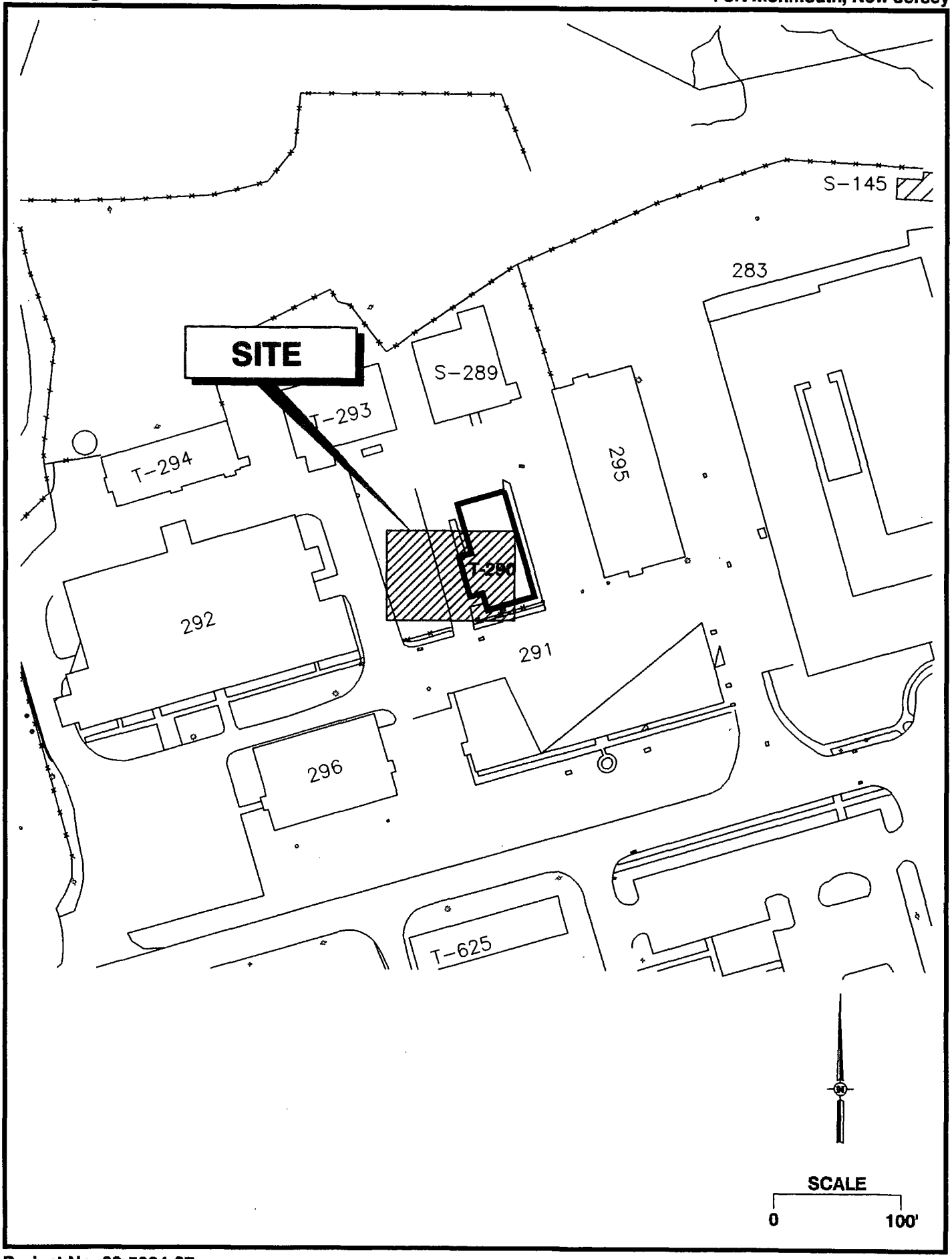

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Figure 2
Building 290

coarse-grained sand that contains abundant rock fragments, minor mica and glauconite (Jablonski). The lower member (Sandy Hook) is a dark gray to black, medium-to-fine grained sand with abundant clay, mica, and glauconite.

The Tinton sand conformably overlies the Red Bank Sand and ranges from a clayey medium to very coarse grained feldspathic quartz and glauconite sand to a glauconitic coarse sand. The color varies from dark yellowish orange or light brown to moderate brown and from light olive to grayish olive. Glauconite may constitute 60 to 80 percent of the sand fraction in the upper part of the unit (Minard, 1969). The upper part of the Tinton is often highly oxidized and iron oxide encrusted (Minard).

Hydrogeology

The water table aquifer in the Main Post area is identified as part of the "composite confining units," or minor aquifers. The minor aquifers include the Navesink formation, Red Bank Sand, Tinton Sand, Hornerstown Sand, Vincentown Formation, Manasquan Formation, Shark River Formation, Piney Point Formation, and the basal clay of the Kirkwood Formation.

Based on records of wells drilled in the Main Post area, water is typically encountered at depths of 2 to 9 feet below ground surface (bgs). According to Jablonski, wells drilled in the Red Bank and Tinton Sands may produce 2 to 25 gallons per minute (gpm). Some well owners have reported acidic water that requires treatment to remove iron.

Due to the proximity of the Atlantic Ocean to Fort Monmouth, shallow groundwater may be tidally influenced and may flow toward creeks and brooks as the tide goes out, and away from creeks and brooks as the tide comes in. However, an abundance of clay lenses and sand deposits were noted in borings installed throughout Fort Monmouth. Therefore the direction of shallow groundwater should be determined on a case by case basis.

1.3 HEALTH AND SAFETY

Before, during, and after all decommissioning activities, hazards at the work site which may have posed a threat to the Health and Safety of all personnel who were involve with, or were affected by, the decommissioning of the UST system were minimized. All areas which posed, or may have been suspected to pose a vapor hazard were monitored by a qualified individual utilizing an organic vapor analyzer (OVA). The individual ascertained if the area was properly vented to render the area safe, as defined by OSHA.

1.4 REMOVAL OF UNDERGROUND STORAGE TANK

1.4.1 General Procedures

- All underground obstructions (utilities, etc.) were marked out by the contractor performing the closure prior to excavation activities.
- All activities were carried out with the greatest regard to safety and health and the safeguarding of the environment.
- All excavated soils were visually examined and screened with an OVA for evidence of contamination. Potentially contaminated soils were identified and logged during closure activities.
- Surface materials (i.e., asphalt, concrete, etc.) were excavated and staged separately from all soil and recycled in accordance with all applicable regulations and laws.
- A Sub-Surface Evaluator from the DPW was present during all site assessment activities.

1.4.2 Underground Storage Tank Excavation and Cleaning

Prior to UST decommissioning activities, surficial soil was removed to expose the UST and associated piping. All free product present in the piping was drained into the UST, and the UST was purged to remove vapors prior to cutting and removal of the piping. After removal of the associated piping, a manway was made in the UST to allow for proper cleaning. The UST was completely emptied of all liquids prior to removal from the ground. Approximately 20 gallons of liquid were transported by Freehold Cartage Inc. to Lionetti Oil Recovery Co. Inc., a NJDEP-approved petroleum recycling and disposal company located in Old Bridge, New Jersey. Refer to Appendix C for the waste manifest (NJA-1907275).

The UST was cleaned prior to removal from the excavation in accordance with the NJDEP-BUST regulations. After the UST was removed from the excavation, it was staged on polyethylene sheeting and examined for holes. No holes or punctures were observed during the inspection by the Sub-Surface Evaluator. Soils surrounding the UST were screened visually and with an OVA for evidence of contamination. Evidence of contamination was observed surrounding the UST.

Soil screening was also performed along the piping associated with the UST. No contamination was noted anywhere along the piping length.

1.5 UNDERGROUND STORAGE TANK TRANSPORTATION AND DISPOSAL

The tank was transported by CUTE Inc. to Monmouth County Reclamation Center for disposal in compliance with all applicable regulations and laws. The UST Disposal Certificate was not available.

The removal contractor labeled the UST prior to transport with the following information:

- site of origin
- contact person
- NJDEP UST Facility ID number
- name of transporter/contact person
- destination site/contact person

1.6 MANAGEMENT OF EXCAVATED SOILS

Based on visual observations, approximately 40 cubic yards of potentially contaminated soil was removed from the excavation on September 12, 1994. On September 13, 1994, approximately 10 additional cubic yards of potentially contaminated soil was removed. All potentially contaminated soils were stockpiled separately from other excavated material and were placed on and covered with polyethylene sheets. Potentially contaminated soils were transported to 490 storage area on Main Post prior to ultimate disposal at Soil Remediation of Philadelphia. Soils that did not exhibit signs of contamination were used as backfill following removal of the UST.

2.0 SITE INVESTIGATION ACTIVITIES

2.1 OVERVIEW

The Site Investigation was managed and carried out by U.S. Army DPW personnel. All TPHC analyses were performed and reported by U.S. Army Fort Monmouth Environmental Laboratory. The VOC, SVOC, and metal analyses were performed and reported by Princeton Laboratory and EMSL Laboratory. All three laboratories are NJDEP-certified testing laboratories. All sampling was performed under the direct supervision of a NJDEP Certified Sub-Surface Evaluator according to the methods described in the NJDEP *Field Sampling Procedures Manual* (1992). Sampling frequency and parameters analyzed complied with he NJDEP-BUST document *Interim Closure Requirements for Underground Storage Tank Systems* (September 1990 and revisions dated November 1, 1991) which was the applicable regulation at the date of the closure. All records of the Site Investigation activities are maintained by the Fort Monmouth DPW Environmental Office.

The following Parties participated in Closure and Site Investigation Activities.

- Closure Contractor: Cleaning Up The Environment Inc. (CUTE Inc.)
  Closure Supervisor: George Bernotsky
  Phone Number: (201)427-2881
  NJDEP Company Certification No.:

- Subsurface Evaluator: Dinkerrai M. Desai
  Employer: U.S. Army, Fort Monmouth
  Phone Number: (908)532-1475
  NJDEP Certification No.: E0002266

- Analytical Laboratory: U.S. Army Fort Monmouth Environmental Laboratory
  Contact Person: Brian K. McKee
  Phone Number: (908)532-4359
  NJDEP Company Certification No.: 13461

- Analytical Laboratory: Princeton Laboratory
  Contact Person: Allan Volk
  Phone Number: (609)452-9050
  NJDEP Company Certification No.: 11118

- Analytical Laboratory: EMSL Analytical, INC.
  Contact Person: Paul Laria
  Phone Number: (908)981-0550
  NJDEP Company Certification No.: 04653

- Hazardous Waste Hauler: Freehold Cartage Inc.
  Contact Person: Barry Olsen
  Phone Number: (908)721-0900
  NJDEP Hazardous Waste Hauler No.: 2265

2.2 FIELD SCREENING/MONITORING

Field screening was performed by a NJDEP Certified Sub-Surface Evaluator using an OVA and visual observations to identify potentially contaminated material. Soil excavated from around the tank and appurtenant piping, as well as the UST excavation sidewalls and bottom, did exhibit evidence of potential contamination. On September 12, 1994, approximately 40 cubic yards of potentially contaminated soil was removed from the excavation due to visual contamination. On September 13, 1994, approximately 10 cubic yards of potentially contaminated soil was removed from the excavation and was stockpiled for disposal.

2.3 SOIL SAMPLING

On September 13, 1994, post-excavation soil samples samples A, B, C, D, E, F, and DUP A were collected from a total of a total of six (6) locations along the sidewalls of the excavation, at a depth of 5.5 feet below ground surface (bgs). Sample H was collected from along the former piping length of the excavation, which was approximately 15 feet in length. The piping samples were collected at a depth of 5.5 feet bgs. All samples were analyzed for total petroleum hydrocarbons (TPHC).

On September 21, 1994, due to elevated TPHC levels, a sample was collected from the north sidewall of the excavation in the vicinity of sample A, and was analyzed for volatile organic compounds plus 15 tentatively identified compounds (VOCs). The sample was collected at a depth of 5.5 feet bgs.

The site assessment was performed by U.S. Army personnel in accordance with the NJDEP *Technical Requirements* and the NJDEP *Field Sampling Procedures Manual.* A summary of sampling activities including parameters analyzed is provided in Table 1. The post-excavation soil samples were collected using polystyrene scoops. Actual soil TPHC values may be higher than reported, due to sample utensil absorbency. If absorbency resulted in reducing the actual soil TPHC concentration by 50 %, the highest soil contaminant would have been 32,400.0 mg/kg, which exceeds the applicable NJDEP soil cleanup standard for total organic contaminants of 10,000 mg/kg. Following soil sampling activities, the samples were chilled and delivered to U.S. Army Fort Monmouth Environmental Laboratory located in Fort Monmouth, New Jersey, for analysis.

2.4 GROUNDWATER SAMPLING

2.4.1 Monitoring Well Installation

One shallow overburden monitoring well (MW-1) was installed at the Building 290 area on July 15, 1994. It was installed approximately 30 feet south off the southeastern corner of Building 290 in the assumed downgradient direction of the former excavation. It was screened in the 2- to- 12.5 foot depth interval, across the water table, which is approximately 3 feet below grade surface.

A second monitoring well (MW-2) was installed at the Building 290 area on August 16, 1994. A monitoring well location map is provided on Figure 4. It was installed approximately 8 feet north off the northwestern corner of Building 290 in the assumed downgradient direction of the former excavation. It was screened in the 1.5- to 11.5- foot depth interval, across the water table, which is approximately 4 feet below grade surface.

The two wells were constructed in accordance with the NJDEP's well construction protocols outlined in its May 1992 *Field Sampling Procedures Manual.* The NJDEP well permits and well construction logs are presented in Appendix D.

Monitoring well MW-1 was constructed with 4-inch (ID) PVC riser and 0.020 slotted PVC well screen. A silica sand pack was installed in the annulus between the borehole wall and the screen. The sand pack was extended approximately 1 foot above the top of the screen. The sand pack above the well screen was graded down to a fine sand to minimize grout intrusion.

The borehole was tremie-grouted with bentonite-cement grout from the top of the sand pack to grade. The well was secured with a water-tight, steel protective casing with a stickup that is approximately 3 feet above ground surface. The steel protective casing was set in place with concrete, which was placed in the remaining open borehole. The elevation of the well riser was surveyed to the nearest 0.01 feet by a New Jersey-licensed surveyor. The well permit number was marked on the well casing as required.

Monitoring well MW-2 was constructed with 4-inch (ID) PVC riser and 0.020 slotted PVC well screen. A silica sand pack was installed in the annulus between the borehole wall and the screen. The sand pack was extended approximately 1 foot above the top of the screen. The sand pack above the well screen was graded down to a fine sand to minimize grout intrusion.

The borehole was tremie-grouted with bentonite-cement grout from the top of the sand pack to grade. The well was secured with a water-tight, flush-mounted locking road box. The locking road box was set in place with concrete, which was placed in the remaining open borehole. The elevation of the well riser was surveyed to the nearest 0.01 feet by a New Jersey-licensed surveyor. The well permit number was marked on the well casing as required.

The monitoring wells were developed using a submersible pump. The wells were pumped for 1 hour or until silt free. All residual soils and liquids generated during monitoring well installation and development program were collected in New Jersey Department of Transportation-approved 55-gallon drums. The drums were placed in a designated secure location for waste characterization and offsite disposal.

2.4.2 Monitoring Well Sampling

Monitoring well MW-1 was sampled on November 8, 1994, November 29, 1994, and December 18, 1995, and analyzed for volatile organic compounds with xylenens (VOCs), tertiary butyl alcohol (TBA), methyl teriary butyl ether (MTBE), and total lead.

Monitoring well MW-2 was sampled on November 18, 1995, and analyzed for volatile organic compounds with xylenes (VOCs), tertiary butyl alcohol, methyl tertiary butyl ether (MTBE), and total lead.

All sampling and analyses were performed in accordance with the NJDEP *Field Sampling Procedures Manual*, and the *Technical Requirements for Site Remediation*, N.J.A.C. 7:26E (*Technical Requirements*).

Prior to sampling each well, the water level was measured to the nearest 0.01 feet, and the distance to the bottom of the well was be measured to the nearest 0.1 feet. The well was checked for floating product (light non-aqueous phase liquids). The well was then purged of three to five well volumes of standing water. Sample volume was then collected using a dedicated decontaminated Teflon bottom-fill bailer attached to PTFE (Teflon)-coated stainless steel.

TABLE 1

SUMMARY OF SAMPLING ACTIVITIES
BUILDING 290, MAIN POST
FORT MONMOUTH, NEW JERSEY

Sample ID	Date of Collection	Matrix	Sample Type	Analytical Parameters (and USEPA Methods) *	Sampling Method
A	9/13/94	Soil	Post-Excavation	TPHC	Polystyrene Scoop
B	9/13/94	Soil	Post-Excavation	TPHC	Polystyrene Scoop
C	9/13/94	Soil	Post-Excavation	TPHC	Polystyrene Scoop
D	9/13/94	Soil	Post-Excavation	TPHC	Polystyrene Scoop
E	9/13/94	Soil	Post-Excavation	TPHC	Polystyrene Scoop
F	9/13/94	Soil	Post-Excavation	TPHC	Polystyrene Scoop
Dup A	9/13/94	Soil	Post-Excavation	TPHC	Polystyrene Scoop
H	9/13/94	Soil	Post-Excavation	TPHC	Polystyrene Scoop
Site A	9/21/94	Soil	Post-Excavation	VOCs	Polystyrene Scoop
290-B-1	3/26/98	Soil	Boring	TPHC	Split Spoon
290-B-2	3/26/98	Soil	Boring	TPHC	Split Spoon
290-B-3	3/26/98	Soil	Boring	TPHC	Split Spoon
290-B-4	3/26/98	Soil	Boring	TPHC	Split Spoon
290-B-5	3/27/98	Soil	Boring	TPHC	Split Spoon
MW-1	11/08/94	Aqueous	Groundwater	Lead, VOCs	Teflon Bottom Bailer
MW-1	11/29/94	Aqueous	Groundwater	Lead, VOCs	Teflon Bottom Bailer
MW-1	12/18/95	Aqueous	Groundwater	VOCs, SVOCs, Metals	Teflon Bottom Bailer
MW-2	11/28/95	Aqueous	Groundwater	VOCs, SVOCs, Metals, Pesticides/PCBs	Teflon Bottom Bailer
MW-1	6/13/97	Aqueous	Groundwater	VOCs, SVOCs, Metals, Pesticides/PCBs	Teflon Bottom Bailer

* Note:

TPHC	Total Petroleum Hydrocarbons (Method 418.1 / soil and aqueous)
VOCs	Volatilte Organic Compounds calibrated for xylenes plus 15 tentatively identified compounds (Method 524.2 / aqueous)
SVOCs	Semivolatile Organic Compounds plus 15 tentatively identified compounds (Method 625 / aqueous)
Metals	(Method SW-846 / aqueous)
Pesticides/PCBs	(Method 608 / aqueous)

3.0 CONCLUSIONS AND RECOMMENDATIONS

3.1 SOIL SAMPLING RESULTS

To evaluate soil conditions following removal of the UST and associated piping, post-excavation soil samples were collected from a total of seven (7) locations on September 13, 1994. All samples were analyzed for TPHC. The post-excavation sampling results were compared to the NJDEP residential direct contact total organic contaminants soil cleanup criteria of 10,000 mg/kg (N.J.A.C. 7:26D and revisions dated February 3, 1994). One sample was collected on September 21, 1994 from the north sidewall of the excavation in the vicinity of sample A and was analyzed for VOCs. The results were compared to the NJDEP Soil Cleanup Criteria. A summary of the analytical results and comparison to the NJDEP soil cleanup criteria is provided in Table 2 and the soil sampling results are shown on Figure 3. The analytical data package is provided in Appendix E.

Post-excavation soil samples (samples C, D, and H) collected from the UST excavation and from below piping associated with the former UST at Building 290 contained TPHC concentrations below the NJDEP residential direct contact total organic contaminants soil cleanup criteria of 10,000 milligrams per kilogram (mg/kg) (N.J.A.C. 7:26D and revisions dated February 3, 1994). Samples C, D, and H, contained levels of TPHC ranging in concentration from 57.3 mg/kg to 730.0 mg/kg. Samples E, and F contained TPHC concentrations of 3,110.0 mg/kg and 9,670.0 mg/kg, respectively. Samples A, B, and DUP A contained levels of TPHC ranging in concentration from 10,400.0 mg/kg to 16,200.0 mg/kg, which exceeded the NJDEP soil cleanup criteria for 10,000 mg/kg. Post-excavation soil sample Site A, collected on September 21, 1994, contained methylene chloride at 0.46 mg/kg, and 2-butanone at 1.5 mg/kg. No other compounds were detected.

3.2 GROUNDWATER SAMPLING RESULTS

The sample collected from MW-1 on November 8, 1994, contained lead at 17.0 micrograms per liter (ug/l). This exceeded the Ground Water Quality Criteria (GWQC) for lead of 10 ug/l. All other groundwater analytical results were either below the detection limit or in compliance with the New Jersey GWQC.

The sample collected on November 29, 1994, contained lead at 3.0 ug/l, which complies with the GWQC for lead. No other compounds were detected.

The sample collected on December 18, 1995, contained methylene chloride at 1.1 ug/l, and cis-1,2-Dichloroethene at 1.1 ug/l. The sample also contained aluminum at 2,200.0 ug/l, barium at 55.0 ug/l, calcium at 4,100.0 ug/l, copper at 60.0 ug/l, iron at 6,400.0 ug/l, potassium at 5,000.0 ug/l, magnesium at 4,500.0 ug/l, manganese at 23.0 ug/l, sodium at 9,200.0 ug/l, and zinc

at 180.0 ug/l. This exceeds the GWQC for alumium of 200.0 ug/l, and iron of 300.0 ug/l. The trip blank and the field blank contained methylene chloride at 1.4 ug/l. No other compounds were detected.

The sample collected from MW-2 on November 28, 1995 contained methylene chloride at 2.3 ug/l, cis-1,2-Dichloroethene at 1.3 ug/l, and chloroform at 0.50 ug/l. The sample also contained silver at 440.0 ug/l, aluminum at 360.0 ug/l, barium at 74.0 ug/l, calcium at 5,300.0 ug/l, iron at 1,100.0 ug/l, magnesium at 4,700.0 ug/l, manganese at 30.0 ug/l, sodium at 11,000.0 ug/l, lead at 1.5 ug/l, and zinc at 100.0 ug/l. This exceeds the GWQC for aluminum of 200.0 ug/l, and iron of 300.0 ug/l. The trip blank and the field blank contained methylene chloride at 0.70 ug/l, and 0.60 ug/l, respectively. The field blank also contained manganese at 32.0 ug/l, lead at 2.8 ug/l, and zinc at 27.0 ug/l. No other compounds were detected.

No product or sheen was observed in MW-1 or MW-2 on any of the sampling dates. The depth to the water table in MW-1 on November 8, 1994 was 8.14 feet below grade, 7.15 feet on November 29, 1994, and 3.06 feet below grade on December 18, 1995. The depth to the water table in MW-2 was 3.18 feet below grade on November 28, 1995.

3.3 CONCLUSIONS AND RECOMMENDATIONS

The analytical results for all post-excavation soil samples collected from the UST closure excavation at Building 290 were below the NJDEP soil cleanup criteria for total organic contaminants.

Based on the post-excavation sampling results, soils with TPHC concentrations exceeding the NJDEP soil cleanup criteria for total organic contaminants of 10,000 mg/kg, do exist in the former location of the UST or associated piping.

The groundwater sample collected on November 8, 1994, contained a lead concentration which exceeded the New Jersey GWQC of 10 ug/l. However, lead was detected at a concentration below the GWQC during a second round of groundwater sampling on November 29, 1994. Based oon the analytical results of the groundwater samples collected on November 29, 1994, November 28, 1995, and December 18, 1995, groundwater quality at the Building 290 UST closure site complies with the New Jersey GWQC.

No further action is proposed in regard to the closure and site assessment of UST No. 081533-64 at Building 290.

TABLE 2

POST-EXCAVATION SOIL SAMPLING RESULTS
BUILDING 290
FT. MONMOUTH, NEW JERSEY

PAGE 1 OF 3

Sample ID/Depth	Sample Laboratory ID	Sample Date	Analysis Date	Compound Name	Sample Quantitation Limit (mg/kg)	Compound of Concern	Result (mg/kg)	NJDEP Soil Cleanup Criteria * (mg/kg)	Exceeds Cleanup Criteria
A/5.5-6.0'	1641.1	9/13/94	9/14/94	Total Solid	--	--	86 %	--	--
				TPHC	130.0	yes	16,200.0	10,000	--
B/5.5-6.0'	1641.2	9/13/94	9/14/94	Total Solid	--	--	88 %	--	--
				TPHC	130.0	yes	11,900.0	10,000	yes
C/5.5-6.0'	1641.3	9/13/94	9/14/94	Total Solid	--	--	83 %	--	--
				TPHC	6.6	yes	730.0	10,000	--
D/5.5-6.0'	1641.4	9/13/94	9/14/94	Total Solid	--	--	82 %	--	--
				TPHC	6.6	yes	126.0	10,000	--
E/5.5-6.0'	1641.5	9/13/94	9/14/94	Total Solid	--	--	84 %	--	--
				TPHC	46.0	yes	3,110.0	10,000	--
F/5.5-6.0'	1641.6	9/13/94	9/14/94	Total Solid	--	--	85 %	--	--
				TPHC	46.0	yes	9,670.0	10,000	--
Dup A/5.5-6.0'	1641.7	9/13/94	9/14/94	Total Solid	--	--	85 %	--	--
				TPHC	46.0	yes	10,400	10,000	yes
H/5.5-6.0'	1641.8	9/13/94	9/14/94	Total Solid	--	--	83 %	--	--
				TPHC	9.9	yes	57.3	10,000	--
290-B-1	3437.02	3/26/98	3/27/98	Total Solid	--	--	92.81 %	--	--
				TPHC	169.0	yes	317.67	10,000	--
290-B-2	3437.03	3/26/98	3/27/98	Total Solid	--	--	84.28 %	--	--
				TPHC	185.0	yes	ND	10,000	--
290-B-3	3437.04	3/26/98	3/27/98	Total Solid	--	--	78.81 %	--	--
				TPHC	196.0	yes	224.45	10,000	--
290-B-4	3437.06	3/26/98	3/27/98	Total Solid	--	--	74.89 %	--	--
				TPHC	196.0	yes	ND	10,000	--
290-B-5	3442.02	3/27/98	3/30/98	Total Solid	--	--	82.86 %	--	--
				TPHC	184.0	yes	ND	10,000	--

Notes:
* Cleanup criteria for total organics
-- Not applicable / does not exceed criteria
ND Not detected above method detection limit
TPHC Total Petroleum Hydrocarbons

Smith Technology Corporation (Project No. 09-5004-08)

TABLE 2

POST-EXCAVATION SOIL SAMPLING RESULTS
BUILDING 290
FT. MONMOUTH, NEW JERSEY
VOLATILE ORGANIC COMPOUNDS

Sample ID/Depth	Sample Date	Analysis Date	Compound Name	Sample Quantitation Limit (mg/kg)	Compound of Concern	Result (mg/kg)	NJDEP Soil Cleanup Criteria * (mg/kg)	Exceeds Cleanup Criteria
Site A/5.5-6.0'	9/21/94	9/23/94	Chloromethane	1.2	--	ND	520/10	--
			Bromomethane	1.2	--	ND	79/1	--
			Vinyl chloride	1.2	--	ND	2/10	--
			Chloroethane	1.2	--	ND	--	--
			Methylene chloride	0.46	--	0.46 J	49/1	--
			Acetone	0.62	--	ND	1,000/100	--
			Carbon Disulfide	0.62	--	ND	--	--
			1,1-Dichloroethene	0.62	--	ND	8/10	
			1,1-Dichloroethane	0.62	--	ND	570/10	--
			1,2-Dichloroethene (total)	0.62	--	ND	79/1	--
			Chloroform	0.62	--	ND	19/1	--
			1,2-Dichloroethane	0.62	--	ND	6/1	--
			2-Butanone	1.5	--	1.5	1,000/50	--
			1,1,1-Trichloroethane	0.62	--	ND	210/50	--
			Carbon Tetrachloride	0.62	--	ND	2/1	--
			Bromodichloromethane	0.62	--	ND	11/1	--
			1,1,2,2-Tetrachloroethane	0.62	--	ND	34/1	--
			1,2-Dichloropropane	0.62	--	ND	10/--	--
			trans-1,3-Dichloropropene	0.62	--	ND	4/1	--
			Trichloroethene	0.62	--	ND	23/1	--
			Dibromochloromethane	0.62	--	ND	110/1	--
			1,1,2-Trichloroethane	0.62	--	ND	22/1	--
			Benzene	0.62	--	ND	3/1	--
			cis-1,3-Dichloropropene	0.62	--	ND	4/1	--
			Bromoform	0.62	--	ND	86/1	--
			2-Hexanone	0.62	--	ND	--	--
			4-Methyl-2-Pentanone	0.62	--	ND	1,000/50	--
			Tetrachloroethene	0.62	--	ND	4/1	--
			Toluene	0.62	--	ND	1,000/500	--
			Chlorobenzene	0.62	--	ND	37/1	--
			Ethylbenzene	0.62	--	ND	1,000/100	--
			Styrene	0.62	--	ND	23/100	--
			Total Xylenes	0.62	--	ND	410/10	--

TABLE 2

POST-EXCAVATION SOIL SAMPLING RESULTS
BUILDING 290
FT. MONMOUTH, NEW JERSEY
VOLATILE ORGANIC COMPOUNDS

Sample ID/Depth	Sample Date	Analysis Date	Compound Name	Sample Quantitation Limit (mg/kg)	Compound of Concern	Result (mg/kg)	NJDEP Soil Cleanup Criteria * (mg/kg)	Exceeds Cleanup Criteria
Field Blank	9/21/94	9/23/94	Chloromethane	0.01	--	ND	520/10	--
			Bromomethane	0.01	--	ND	79/1	--
			Vinyl chloride	0.01	--	ND	2/10	--
			Chloroethane	0.01	--	ND	--	--
			Methylene chloride	0.005	--	ND	49/1	--
			Acetone	0.005	--	ND	1,000/100	--
			Carbon Disulfide	0.005	--	ND	--	--
			1,1-Dichloroethene	0.005	--	ND	8/10	
			1,1-Dichloroethane	0.005	--	ND	570/10	--
			1,2-Dichloroethene (total)	0.005	--	ND	79/1	--
			Chloroform	0.005	--	ND	19/1	--
			1,2-Dichloroethane	0.005	--	ND	6/1	--
			2-Butanone	0.005	--	ND	1,000/50	--
			1,1,1-Trichloroethane	0.005	--	ND	210/50	--
			Carbon Tetrachloride	0.005	--	ND	2/1	--
			Bromodichloromethane	0.005	--	ND	11/1	--
			1,1,2,2-Tetrachloroethane	0.005	--	ND	34/1	--
			1,2-Dichloropropane	0.005	--	ND	10/--	--
			trans-1,3-Dichloropropene	0.005	--	ND	4/1	--
			Trichloroethene	0.005	--	ND	23/1	--
			Dibromochloromethane	0.005	--	ND	110/1	--
			1,1,2-Trichloroethane	0.005	--	ND	22/1	--
			Benzene	0.005	--	ND	3/1	--
			cis-1,3-Dichloropropene	0.005	--	ND	4/1	--
			Bromoform	0.005	--	ND	86/1	--
			2-Hexanone	0.005	--	ND	--	--
			4-Methyl-2-Pentanone	0.005	--	ND	1,000/50	--
			Tetrachloroethene	0.005	--	ND	4/1	--
			Toluene	0.005	--	ND	1,000/500	--
			Chlorobenzene	0.005	--	ND	37/1	--
			Ethylbenzene	0.005	--	ND	1,000/100	--
			Styrene	0.005	--	ND	23/100	--
			Total Xylenes	0.005	--	ND	410/10	--

Notes:
* Residential Direct Contact / Impact to Groundwater
-- Not applicable / does not exceed criteria
(ND) Indicates compound is not detected

TABLE 3

GROUNDWATER SAMPLING RESULTS
BUILDING 290, MAIN POST, MW-1
FORT MONMOUTH, NEW JERSEY

Sample ID/Depth	Sample Date	Analysis Date	Compound Name	Sample Quantitation Limit (ug/l)	Compound of Concern	Result (ug/l)	GWQC (ug/l)	Exceeds Cleanup Criteria
MW-1	6/13/97	6/19/97	Lead	17.0	--	17.0	10	YES
			VOLATILE ORGANICS:					
			Chloromethane	ND	--	ND	--	--
			Bromomethane	ND	--	ND	--	--
			Vinyl chloride	ND	--	ND	5	--
			Chloroethane	ND	--	ND	--	--
			Methylene chloride	ND	--	ND	2	--
			1,1-Dichloroethene	ND	--	ND	2	--
			1,1-Dichloroethane	ND	--	ND	70	--
			Chloroform	ND	--	ND	6	--
			1,2-Dichloroethane	ND	--	ND	2	--
			1,1,1-Trichloroethane	ND	--	ND	30	--
			Carbon Tetrachloride	ND	--	ND	2	--
			Bromodichloromethane	ND	--	ND	1	--
			1,2-Dichloropropane	ND	--	ND	1	--
			cis-1,3-Dichloropropene	ND	--	ND	NA	--
			Trichloroethene	ND	--	ND	1	--
			Dibromochloromethane	ND	--	ND	10	--
			1,1,2-Trichloroethane	ND	--	ND	3	--
			Benzene	ND	--	ND	1	--
			trans-1,3-Dichloropropene	ND	--	ND	NA	--
			Bromoform	ND	--	ND	4	--
			Tetrachloroethene	ND	--	ND	1	--
			1,1,2,2-Tetrachloroethane	ND	--	ND	2	--
			Toluene	ND	--	ND	1,000	--
			Chlorobenzene	ND	--	ND	4	--
			Ethylbenzene	ND	--	ND	700	--
			Xylene (total)	ND	--	ND	40	--
			Trichloromonofluoromethane	ND	--	ND	--	--
			Acrolein	ND	--	ND	NA	--
			Acrylonitrile	ND	--	ND	50	--
			Tertiary Butyl Alcohol	ND	--	ND	--	--
			Methyl Tertiary Butyl Ether	ND	--	ND	--	--
			1,3-Dichlorobenzene	ND	--	ND	600	--
			1,4-Dichlorobenzene	ND	--	ND	75	--
			1,2-Dichlorobenzene	ND	--	ND	600	--
			2-Chloroethylvinyl Ether	ND	--	ND	--	--
			Trans,1,2-Dichloroethene	ND	--	ND	100	--
			VOLATILE TICS:					
			Unknown Siloxane	--	--	29 J	--	--
			Unknown Siloxane	--	--	110 J	--	--
			Unknown Siloxane	--	--	52 J	--	--
			TOTOAL TICS:	--	--	191	--	--

TABLE 3

GROUNDWATER SAMPLING RESULTS
BUILDING 290, MAIN POST, TRIP BLANK
FORT MONMOUTH, NEW JERSEY

Sample ID/Depth	Sample Date	Analysis Date	Compound Name	Sample Quantitation Limit (ug/l)	Compound of Concern	Result (ug/l)	GWQC (ug/l)	Exceeds Cleanup Criteria
TRIP BLANK	6/13/97	6/19/97	Lead	--	--	NT	10	--
			VOLATILE ORGANICS:					
			Chloromethane	ND	--	ND	--	--
			Bromomethane	ND	--	ND	--	--
			Vinyl chloride	ND	--	ND	5	--
			Chloroethane	ND	--	ND	--	--
			Methylene chloride	ND	--	ND	2	--
			1,1-Dichloroethene	ND	--	ND	2	--
			1,1-Dichloroethane	ND	--	ND	70	--
			Chloroform	ND	--	ND	6	--
			1,2-Dichloroethane	ND	--	ND	2	--
			1,1,1-Trichloroethane	ND	--	ND	30	--
			Carbon Tetrachloride	ND	--	ND	2	--
			Bromodichloromethane	ND	--	ND	1	--
			1,2-Dichloropropane	ND	--	ND	1	--
			cis-1,3-Dichloropropene	ND	--	ND	NA	--
			Trichloroethene	ND	--	ND	1	--
			Dibromochloromethane	ND	--	ND	10	--
			1,1,2-Trichloroethane	ND	--	ND	3	--
			Benzene	ND	--	ND	1	--
			trans-1,3-Dichloropropene	ND	--	ND	NA	--
			Bromoform	ND	--	ND	4	--
			Tetrachloroethene	ND	--	ND	1	--
			1,1,2,2-Tetrachloroethane	ND	--	ND	2	--
			Toluene	ND	--	ND	1,000	--
			Chlorobenzene	ND	--	ND	4	--
			Ethylbenzene	ND	--	ND	700	--
			Xylene (total)	ND	--	ND	40	--
			Trichloromonofluoromethane	ND	--	ND	--	--
			Acrolein	ND	--	ND	NA	--
			Acrylonitrile	ND	--	ND	50	--
			Tertiary Butyl Alcohol	ND	--	ND	--	--
			Methyl Tertiary Butyl Ether	ND	--	ND	--	--
			1,3-Dichlorobenzene	ND	--	ND	600	--
			1,4-Dichlorobenzene	ND	--	ND	75	--
			1,2-Dichlorobenzene	ND	--	ND	600	--
			2-Chloroethylvinyl Ether	ND	--	ND	--	--
			Trans,1,2-Dichloroethene	ND	--	ND	100	--
			VOLATILE TICS:					
			NONE FOUND	--	--	--	--	--

TABLE 3

GROUNDWATER SAMPLING RESULTS
BUILDING 290, MAIN POST, FIELD BLANK
FORT MONMOUTH, NEW JERSEY

Sample ID/Depth	Sample Date	Analysis Date	Compound Name	Sample Quantitation Limit (ug/l)	Compound of Concern	Result (ug/l)	GWQC (ug/l)	Exceeds Cleanup Criteria
FIELD BLANK	6/13/97	6/19/97	Lead	4.1	--	4.1 B	10	--
			VOLATILE ORGANICS:					
			Chloromethane	ND	--	ND	--	--
			Bromomethane	ND	--	ND	--	--
			Vinyl chloride	ND	--	ND	5	--
			Chloroethane	ND	--	ND	--	--
			Methylene chloride	ND	--	4	2	YES
			1,1-Dichloroethene	ND	--	ND	2	--
			1,1-Dichloroethane	ND	--	ND	70	--
			Chloroform	ND	--	ND	6	--
			1,2-Dichloroethane	ND	--	ND	2	--
			1,1,1-Trichloroethane	ND	--	ND	30	--
			Carbon Tetrachloride	ND	--	ND	2	--
			Bromodichloromethane	ND	--	ND	1	--
			1,2-Dichloropropane	ND	--	ND	1	--
			cis-1,3-Dichloropropene	ND	--	ND	NA	--
			Trichloroethene	ND	--	ND	1	--
			Dibromochloromethane	ND	--	ND	10	--
			1,1,2-Trichloroethane	ND	--	ND	3	--
			Benzene	ND	--	ND	1	--
			trans-1,3-Dichloropropene	ND	--	ND	NA	--
			Bromoform	ND	--	ND	4	--
			Tetrachloroethene	ND	--	ND	1	--
			1,1,2,2-Tetrachloroethane	ND	--	ND	2	--
			Toluene	ND	--	ND	1,000	--
			Chlorobenzene	ND	--	ND	4	--
			Ethylbenzene	ND	--	ND	700	--
			Xylene (total)	ND	--	ND	40	--
			Trichloromonofluoromethane	ND	--	ND	--	--
			Acrolein	ND	--	ND	NA	--
			Acrylonitrile	ND	--	ND	50	--
			Tertiary Butyl Alcohol	ND	--	ND	--	--
			Methyl Tertiary Butyl Ether	ND	--	ND	--	--
			1,3-Dichlorobenzene	ND	--	ND	600	--
			1,4-Dichlorobenzene	ND	--	ND	75	--
			1,2-Dichlorobenzene	ND	--	ND	600	--
			2-Chloroethylvinyl Ether	ND	--	ND	--	--
			Trans,1,2-Dichloroethene	ND	--	ND	100	--
			VOLATILE TICS:					
			Unknown Hydrocarbon	--	--	3	--	--

TABLE 3

GROUNDWATER SAMPLING RESULTS
BUILDING 290, MAIN POST, MW-1
FORT MONMOUTH, NEW JERSEY

Sample ID/Depth	Sample Date	Analysis Date	Compound Name	Sample Quantitation Limit (ug/l)	Compound of Concern	Result (ug/l)	GWQC (ug/l)	Exceeds Cleanup Criteria
MW-1	6/13/97	6/19/97	Lead	3.0	--	ND	10	--
			VOLATILE ORGANICS:					
			Chloromethane	2	--	ND	--	--
			Bromomethane	1	--	ND	--	--
			Vinyl chloride	1	--	ND	5	--
			Chloroethane	1	--	ND	--	--
			Methylene chloride	3	--	ND	2	--
			1,1-Dichloroethene	ND	--	ND	2	--
			1,1-Dichloroethane	ND	--	ND	70	--
			Chloroform	ND	--	ND	6	--
			1,2-Dichloroethane	ND	--	ND	2	--
			1,1,1-Trichloroethane	ND	--	ND	30	--
			Carbon Tetrachloride	ND	--	ND	2	--
			Bromodichloromethane	ND	--	ND	1	--
			1,2-Dichloropropane	ND	--	ND	1	--
			cis-1,3-Dichloropropene	ND	--	ND	NA	--
			Trichloroethene	ND	--	ND	1	--
			Dibromochloromethane	ND	--	ND	10	--
			1,1,2-Trichloroethane	ND	--	ND	3	--
			Benzene	ND	--	ND	1	--
			trans-1,3-Dichloropropene	ND	--	ND	NA	--
			Bromoform	ND	--	ND	4	--
			Tetrachloroethene	ND	--	ND	1	--
			1,1,2,2-Tetrachloroethane	ND	--	ND	2	--
			Toluene	ND	--	ND	1,000	--
			Chlorobenzene	ND	--	ND	4	--
			Ethylbenzene	ND	--	ND	700	--
			Xylene (total)	ND	--	ND	40	--
			Trichloromonofluoromethane	ND	--	ND	--	--
			Acrolein	ND	--	ND	NA	--
			Acrylonitrile	ND	--	ND	50	--
			Tertiary Butyl Alcohol	ND	--	ND	--	--
			Methyl Tertiary Butyl Ether	ND	--	ND	--	--
			1,3-Dichlorobenzene	ND	--	ND	600	--
			1,4-Dichlorobenzene	ND	--	ND	75	--
			1,2-Dichlorobenzene	ND	--	ND	600	--
			2-Chloroethylvinyl Ether	ND	--	ND	--	--
			Trans,1,2-Dichloroethene	ND	--	ND	100	--
			VOLATILE TICS:					
			Unknown	--	--	9 J	--	--

TABLE 3

GROUNDWATER SAMPLING RESULTS
BUILDING 290, MAIN POST, TRIP BLANK
FORT MONMOUTH, NEW JERSEY

Sample ID/Depth	Sample Date	Analysis Date	Compound Name	Sample Quantitation Limit (ug/l)	Compound of Concern	Result (ug/l)	GWQC (ug/l)	Exceeds Cleanup Criteria
Trip Blank	11/29/94	12/01/94	VOLATILE ORGANICS:					
			Chloromethane	ND	--	ND	--	--
			Bromomethane	ND	--	ND	--	--
			Vinyl chloride	ND	--	ND	5	--
			Chloroethane	ND	--	ND	--	--
			Methylene chloride	ND	--	3.0	2	YES
			1,1-Dichloroethene	ND	--	ND	2	--
			1,1-Dichloroethane	ND	--	ND	70	--
			Chloroform	ND	--	ND	6	--
			1,2-Dichloroethane	ND	--	ND	2	--
			1,1,1-Trichloroethane	ND	--	ND	30	--
			Carbon Tetrachloride	ND	--	ND	2	--
			Bromodichloromethane	ND	--	ND	1	--
			1,2-Dichloropropane	ND	--	ND	1	--
			cis-1,3-Dichloropropene	ND	--	ND	NA	--
			Trichloroethene	ND	--	ND	1	--
			Dibromochloromethane	ND	--	ND	10	--
			1,1,2-Trichloroethane	ND	--	ND	3	--
			Benzene	ND	--	ND	1	--
			trans-1,3-Dichloropropene	ND	--	ND	NA	--
			Bromoform	ND	--	ND	4	--
			Tetrachloroethene	ND	--	ND	1	--
			1,1,2,2-Tetrachloroethane	ND	--	ND	2	--
			Toluene	ND	--	ND	1,000	--
			Chlorobenzene	ND	--	ND	4	--
			Ethylbenzene	ND	--	ND	700	--
			Xylene (total)	ND	--	ND	40	--
			Trichloromonofluoromethane	ND	--	ND	--	--
			Acrolein	ND	--	ND	NA	--
			Acrylonitrile	ND	--	ND	50	--
			Tertiary Butyl Alcohol	ND	--	ND	--	--
			Methyl Tertiary Butyl Ether	ND	--	ND	--	--
			1,3-Dichlorobenzene	ND	--	ND	600	--
			1,4-Dichlorobenzene	ND	--	ND	75	--
			1,2-Dichlorobenzene	ND	--	ND	600	--
			2-Chloroethylvinyl Ether	ND	--	ND	--	--
			Trans,1,2-Dichloroethene	ND	--	ND	100	--
			VOLATILE TICS:					
			Unknown Hydrocarbon	--	--	13.0 J	--	--

TABLE 3

GROUNDWATER SAMPLING RESULTS
BUILDING 290, MAIN POST, FIELD BLANK
FORT MONMOUTH, NEW JERSEY

Sample ID/Depth	Sample Date	Analysis Date	Compound Name	Sample Quantitation Limit (ug/l)	Compound of Concern	Result (ug/l)	GWQC (ug/l)	Exceeds Cleanup Criteria
Field Blank	11/29/94	12/01/94	Lead	3.0	--	ND	10	--
			VOLATILE ORGANICS:					
			Chloromethane	ND	--	ND	--	--
			Bromomethane	ND	--	ND	--	--
			Vinyl chloride	ND	--	ND	5	--
			Chloroethane	ND	--	ND	--	--
			Methylene chloride	ND	--	ND	2	--
			1,1-Dichloroethene	ND	--	ND	2	--
			1,1-Dichloroethane	ND	--	ND	70	--
			Chloroform	ND	--	ND	6	--
			1,2-Dichloroethane	ND	--	ND	2	--
			1,1,1-Trichloroethane	ND	--	ND	30	--
			Carbon Tetrachloride	ND	--	ND	2	--
			Bromodichloromethane	ND	--	ND	1	--
			1,2-Dichloropropane	ND	--	ND	1	--
			cis-1,3-Dichloropropene	ND	--	ND	NA	--
			Trichloroethene	ND	--	ND	1	--
			Dibromochloromethane	ND	--	ND	10	--
			1,1,2-Trichloroethane	ND	--	ND	3	--
			Benzene	ND	--	ND	1	--
			trans-1,3-Dichloropropene	ND	--	ND	NA	--
			Bromoform	ND	--	ND	4	--
			Tetrachloroethene	ND	--	ND	1	--
			1,1,2,2-Tetrachloroethane	ND	--	ND	2	--
			Toluene	ND	--	ND	1,000	--
			Chlorobenzene	ND	--	ND	4	--
			Ethylbenzene	ND	--	ND	700	--
			Xylene (total)	ND	--	ND	40	--
			Trichloromonofluoromethane	ND	--	ND	--	--
			Acrolein	ND	--	ND	NA	--
			Acrylonitrile	ND	--	ND	50	--
			Tertiary Butyl Alcohol	ND	--	ND	--	--
			Methyl Tertiary Butyl Ether	ND	--	ND	--	--
			1,3-Dichlorobenzene	ND	--	ND	600	--
			1,4-Dichlorobenzene	ND	--	ND	75	--
			1,2-Dichlorobenzene	ND	--	ND	600	--
			2-Chloroethylvinyl Ether	ND	--	ND	--	--
			Trans,1,2-Dichloroethene	ND	--	ND	100	--
			VOLATILE TICS:					
			Unknown Hydrocarbon	--	--	3.0 J	--	--
			Unknown Hydrocarbon	--	--	12.0 J	--	--
			TOTAL TICS:	--	--	15.0	--	--

TABLE 3

GROUNDWATER SAMPLING RESULTS
BUILDING 290, MAIN POST, MW-2
FT. MONMOUTH, NEW JERSEY

Sample ID	Sample Date	Analysis Date	Compound Name	Sample Quantitation Limit (ug/l)	Compound of Concern	Result (ug/l)	GWQC (ug/l)	Exceeds Criteria
MW-2	11/28/95	12/11/95	VOLATILE ORGANICS:					
			Dichlorodifluoromethane	0.50	--	ND	--	--
			Chloromethane	0.50	--	ND	--	--
			Vinyl Chloride	0.50	--	ND	--	--
			Bromomethane	0.50	--	ND	5	--
			Chloroethane	0.50	--	ND	--	--
			Trichlorofluoromethane	0.50	--	ND	--	--
			1,1-Dichloroethene	0.50	--	ND	2	--
			Methylene Chloride	2.3	--	2.3 B	100	--
			1,2-Dichloroethene (trans)	0.50	--	ND	2	--
			1,1 Dichloroethane	0.50	--	ND	70	--
			2,2-Dichloropropane	0.50	--	ND	--	--
			cis-1,2-Dichloroethene	1.3	--	1.3	--	--
			Bromochloromethane	0.50	--	ND	10	--
			Chloroform	0.50	--	ND	6	--
			1,1,1-Trichloroethane	0.50	--	ND	--	--
			Carbon Tetrachloride	0.50	--	ND	2	--
			1,1-Dichloropropene	0.50	--	ND	30	--
			Benzene	0.50	--	ND	--	--
			1,2-Dichloroethane	0.50	--	ND	2	--
			Trichloroethene	0.50	--	ND	1	--
			1,2-Dichloropropane	0.50	--	ND	1	--
			Dibromomethane	0.50	--	ND	NA	--
			Bromodichloromethane	0.50	--	ND	--	--
			cis-1,3-Dichloropropene	0.50	--	ND	1	--
			Toluene	0.50	--	ND	10	--
			trans-1,3-Dichloropropene	0.50	--	ND	3	--
			1,1,2-Trichloroethane	0.50	--	ND	1	--
			Tetrachloroethene	0.50	--	ND	NA	--
			1,3-Dichloropropane	0.50	--	ND	4	--
			Dibromochloromethane	0.50	--	ND	10	--
			1,2-Dibromomethane	0.50	--	ND	1	--
			Chlorobenzene	0.50	--	ND	2	--
			1,1,2,2-Tetrachloroethane	0.50	--	ND	1,000	--
			Ethylbenzene	0.50	--	ND	--	--
			Xylenes (Total)	0.50	--	ND	4	--
			Styrene	0.50	--	ND	700	--
			Bromoform	0.50	--	ND	40	--
			Isopropylbenzene	0.50	--	ND	100	--
			Bromobenzene	0.50	--	ND	--	--
			1,1,2,2-Tetrachloroethane	0.50	--	ND	--	--
			1,2,3-Trichloropropane	0.50	--	ND	--	--
			n-Propylbenzene	0.50	--	ND	--	--
			2-Chlorotoluene	0.50	--	ND	--	--

TABLE 3

GROUNDWATER SAMPLING RESULTS
BUILDING 290, MAIN POST, MW-2
FORT MONMOUTH, NEW JERSEY
VOLATILE ORGANICS (Continued)

Sample ID	Sample Date	Analysis Date	Compound Name	Sample Quantitation Limit (ug/l)	Compound of Concern	Result (ug/l)	GWQC (ug/l)	Exceeds Criteria
MW-2	11/28/95	12/11/95	VOLATILES (CONTINUED)					
			4-Chlorotoluene	0.50	--	ND	--	--
			1,3,5-Trimethylbenzene	0.50	--	ND	--	--
			tert-Butylbenzene	0.50	--	ND	--	--
			1,2,4-Trimethylbenzene	0.50	--	ND	--	--
			sec-Butylbenzene	0.50	--	ND	--	--
			1,3-Dichlorobenzene	0.50	--	ND	600	--
			4-Isopropyltoluene	0.50	--	ND	75	--
			1,4-Dichlorobenzene	0.50	--	ND	--	--
			1,2-Dichlorobenzene	0.50	--	ND	600	--
			N-Butylbenzene	0.50	--	ND	--	--
			1,2-Dibromo-3-chloropropane	0.50	--	ND	NA	--
			1,2,4-Trichlorobenzene	0.50	--	ND	9	--
			Hexachlorobutadiene	0.50	--	ND	1	--
			Naphthalene	0.50	--	ND	--	--
			1,2,3-Trichlorobenzene	0.50	--	ND	--	--
			Methy-tertiary butyl ether	0.50	--	ND	--	--
			tertiary-Butyl alcohol	2.0	--	ND	--	--
			VOLATILE TICS:					
			Column Bleed	--	--	1 J	--	--
			Column Bleed	--	--	1 J	--	--
			TOTAL TICS:	--	--	2	--	--

TABLE 3

GROUNDWATER SAMPLING RESULTS
BUILDING 290, MAIN POST, MW-2
FORT MONMOUTH, NEW JERSEY

Sample ID	Sample Date	Analysis Date	Compound Name	Sample Quantitation Limit (ug/l)	Compound of Concern	Result (ug/l)	GWQC (ug/l)	Exceeds Criteria
MW-2	11/28/95	12/11/95	SEMIVOLATILES:					
			Phenol	10	--	ND	4000	--
			bis(2 chloroethyl)Ether	10	--	ND	20	--
			2-Chlorophenol	10	--	ND	40	--
			1,3-Dichlorobenzene	10	--	ND	600	--
			1,4-Dichlorobenzene	10	--	ND	75	--
			1,2-Dichlorobenzene	10	--	ND	600	--
			2-Methylphenol	10	--	ND	--	--
			bis(2-chloroisopropyl)Ether	10	--	ND	300	--
			4-Methylphenol	10	-	ND	--	--
			N-Nitroso-Di-n-propylamine	10	--	ND	20	--
			Hexachloroethane	10	--	ND	10	--
			Nitrobenzene	10	--	ND	10	--
			Isophorone	10	--	ND	100	--
			2-Nitrophenol	10	--	ND	--	--
			2,4-Dimethylphenol	10	--	ND	100	--
			bis(2-Chloroethoxy)Methane	10	--	ND	--	--
			2,4-Dichlorophenol	10	--	ND	20	--
			1,2,4-Trichlorobenzene	10	--	ND	9	--
			Naphthalene	10	--	ND	--	--
			4-Chloroaniline	10	--	ND	--	--
			Hexachlorobutadiene	10	--	ND	1	--
			4-Chloro-3-methylphenol	10	--	ND	--	--
			2-methylnaphthanene	10	--	ND	--	--
			Hexachlorocyclopentadiene	10	--	ND	50	--
			2,4,6-Trichlorophenol	10	--	ND	20	--
			2,4,5-Trichlorophenol	25	--	ND	700	--
			2-Chloronaphthalene	10	--	ND	--	--
			2-Nitroanline	25	--	ND	--	--
			Dimethyl Phthalate	10	--	ND	--	--
			Acenaphthylene	10	--	ND	NA	--
			2,6-Dinitrotoluene	10	--	ND	NA	--
			3-Nitroaniline	25	--	ND	--	--
			Acenaphthene	10	--	ND	400	--
			2,4-Dinitrophenol	25	--	ND	40	--
			4-Nitrophenol	25	--	ND	--	--
			Dibenzofuran	10	--	ND	--	--
			2,4-Dinitrotoluene	10	--	ND	10	--
			Diethylphthalate	10	--	ND	5,000	--
			Fluorene	10	--	ND	300	--
			4-Chlorophenyl-phenlyether	10	--	ND	--	--
			4-Nitroaniline	25	--	ND	--	--
			4,6-Dinitro-2-methylphenol	25	--	ND	--	--

TABLE 3

GROUNDWATER SAMPLING RESULTS
BUILDING 290, MAIN POST, MW-2
FORT MONMOUTH, NEW JERSEY

Sample ID	Sample Date	Analysis Date	Compound Name	Sample Quantitation Limit (ug/l)	Compound of Concern	Result (ug/l)	GWQC (ug/l)	Exceeds Criteria
MW-2	11/28/95	12/11/95	SEMIVOLATILES (Continued)					
			N-Nitrosodiphenylamine	10	--	ND	20	--
			4-Bromophenyl-phenylether	10	--	ND	--	--
			Hexachlorobenzene	10	--	ND	10	--
			Pentachlorophenol	25	--	ND	1	--
			Phenanthrene	10	--	ND	NA	--
			Anthracene	10	--	ND	2,000	--
			Carbazole	10	--	ND	--	--
			Di-n-butylphthalate	10	--	ND	900	--
			Fluoranthene	10	--	ND	300	--
			Pyrene	10	--	ND	200	--
			Butylbenzylphthalate	10	--	ND	100	--
			Benzo(a)Anthracene	10	--	ND	NA	--
			3,3-Dichlorobenzidine	20	--	ND	60	--
			Chrysene	10	--	ND	NA	--
			bis(2-Ethylhexyl)Phthalate	10	--	ND	30	--
			Di-n-Octyl Phthalate	10	--	ND	100	--
			Benzo(b)Fluoranthene	10	--	ND	NA	--
			Benzo(k)Fluoranthene	10	--	ND	NA	--
			Benzo(a)Pyrene	10	--	ND	NA	--
			Indeno(1,2,3-cd)pyrene	10	--	ND	NA	--
			Dibenzo(a,h)anthracene	10	--	ND	NA	--
			Benzo(g,h,i)perylene	10	--	ND	NA	--
			SEMIVOLATILE TICS:					
			Hexadeconoic acid	--	--	28 J	--	--
			Unknown Hydrocarbon	--	--	4 J	--	--
			TOTAL TICS:	--	--	32	--	--

TABLE 3

GROUNDWATER SAMPLING RESULTS
BUILDING 290, MAIN POST, MW-2
FORT MONMOUTH, NEW JERSEY

Sample ID	Sample Date	Analysis Date	Compound Name	Sample Quantitation Limit (ug/l)	Compound of Concern	Result (ug/l)	GWQC (ug/l)	Exceeds Criteria
MW-2	11/28/95	12/07/95	METALS					
			Silver	440	--	440	NA	--
			Aluminum	360	--	360	200	YES
			Arsenic	5	--	ND	8	--
			Barium	74	--	74	2,000	--
			Beryllium	5	--	ND	20	--
			Calcium	5,300	--	5,300	--	--
			Cadmium	10	--	ND	4	--
			Cobalt	50	--	ND	--	--
			Chromium	50	--	ND	100	--
			Copper	50	--	ND	1,000	--
			Iron	1,100	--	1,100	300	YES
			Mercury	1	--	ND	2	--
			Potassium	3,000	--	ND	--	--
			Magnesium	4,700	--	4,700	--	--
			Manganese	30	--	30	50	--
			Sodium	11,000	--	11,000	50,000	--
			Nickel	50	--	ND	100	--
			Lead	1.5	--	1.5	10	--
			Antimony	5	--	ND	20	--
			Selenium	5	--	ND	50	--
			Thallium	2	--	ND	10	--
			Vanadium	50	--	ND	--	--
			Zinc	100	--	100	5,000	--
			Cyanide (Total)	10	--	ND	200	--

TABLE 3

GROUNDWATER SAMPLING RESULTS
BUILDING 290, MAIN POST, MW-2
FORT MONMOUTH, NEW JERSEY

Sample ID	Sample Date	Analysis Date	Compound Name	Sample Quantitation Limit (ug/l)	Compound of Concern	Result (ug/l)	GWQC (ug/l)	Exceeds Criteria
MW-2	11/28/95	12/07/95	PESTICIDES/PCBs					
			alpha-BHC	0.02	--	ND	0.02	--
			beta-BHC	0.04	--	ND	0.2	--
			delta-BHC	0.02	--	ND	--	--
			gamma-BHC (Lindane)	0.03	--	ND	0.2	--
			Heptachlor	0.02	--	ND	0.4	--
			Aldrin	0.04	--	ND	0.04	--
			Heptachlor epoxide	0.05	--	ND	0.2	--
			Endosulfan I	0.02	--	ND	0.4	--
			Dieldrin	0.03	--	ND	0.03	--
			4,4'-DDE	0.04	--	ND	0.1	--
			Endrin	0.04	--	ND	2.0	--
			Endosulfan II	0.04	--	ND	0.4	--
			4,4'-DDD	0.04	--	ND	0.1	--
			Endosulfan sulfate	0.08	--	ND	0.4	--
			4,4'-DDT	0.04	--	ND	0.1	--
			Endrin aldehyde	0.1	--	ND	--	--
			Chlordane	0.1	--	ND	0.5	--
			Toxaphene	1.0	--	ND	3.0	--
			Aroclor-1016	1.0	--	ND	--	--
			Aroclor-1221	1.0	--	ND	--	--
			Aroclor-1232	1.0	--	ND	--	--
			Aroclor-1242	1.0	--	ND	--	--
			Aroclor-1248	1.0	--	ND	--	--
			Aroclor-1254	1.0	--	ND	--	--
			Aroclor-1260	1.0	--	ND	--	--

TABLE 3

GROUNDWATER SAMPLING RESULTS
BUILDING 290, MAIN POST, TRIP BLANK
FT. MONMOUTH, NEW JERSEY

Sample ID	Sample Date	Analysis Date	Compound Name	Sample Quantitation Limit (ug/l)	Compound of Concern	Result (ug/l)	GWQC (ug/l)	Exceeds Criteria
Trip Blank	11/28/95	12/11/95	VOLATILE ORGANICS					
			Dichlorodifluoromethane	0.50	--	ND	--	--
			Chloromethane	0.50	--	ND	--	--
			Vinyl Chloride	0.50	--	ND	--	--
			Bromomethane	0.50	--	ND	5	--
			Chloroethane	0.50	--	ND	--	--
			Trichlorofluoromethane	0.50	--	ND	--	--
			1,1-Dichloroethene	0.50	--	ND	2	--
			Methylene Chloride	3.0	--	3.0 B	100	--
			1,2-Dichloroethene (trans)	0.50	--	ND	2	--
			1,1 Dichloroethane	0.50	--	ND	70	--
			2,2-Dichloropropane	0.50	--	ND	--	--
			cis-1,2-Dichloroethene	0.50	--	ND	--	--
			Bromochloromethane	0.50	--	ND	10	--
			Chloroform	0.50	--	ND	6	--
			1,1,1-Trichloroethane	0.50	--	ND	--	--
			Carbon Tetrachloride	0.50	--	ND	2	--
			1,1-Dichloropropene	0.50	--	ND	30	--
			Benzene	0.50	--	ND	--	--
			1,2-Dichloroethane	0.50	--	ND	2	--
			Trichloroethene	0.50	--	ND	1	--
			1,2-Dichloropropane	0.50	--	ND	1	--
			Dibromomethane	0.50	--	ND	NA	--
			Bromodichloromethane	0.50	--	ND	--	--
			cis-1,3-Dichloropropene	0.50	--	ND	1	--
			Toluene	0.50	--	ND	10	--
			trans-1,3-Dichloropropene	0.50	--	ND	3	--
			1,1,2-Trichloroethane	0.50	--	ND	1	--
			Tetrachloroethene	0.50	--	ND	NA	--
			1,3-Dichloropropane	0.50	--	ND	4	--
			Dibromochloromethane	0.50	--	ND	10	--
			1,2-Dibromomethane	0.50	--	ND	1	--
			Chlorobenzene	0.50	--	ND	2	--
			1,1,2,2-Tetrachloroethane	0.50	--	ND	1,000	--
			Ethylbenzene	0.50	--	ND	--	--
			Xylenes (Total)	0.50	--	ND	4	--
			Styrene	0.50	--	ND	700	--
			Bromoform	0.50	--	ND	40	--
			Isopropylbenzene	0.50	--	ND	100	--
			Bromobenzene	0.50	--	ND	--	--
			1,1,2,2-Tetrachloroethane	0.50	--	ND	--	--
			1,2,3-Trichloropropane	0.50	--	ND	--	--
			n-Propylbenzene	0.50	--	ND	--	--
			2-Chlorotoluene	0.50	--	ND	--	--

TABLE 3

GROUNDWATER SAMPLING RESULTS
BUILDING 290, MAIN POST, TRIP BLANK
FORT MONMOUTH, NEW JERSEY

Sample ID	Sample Date	Analysis Date	Compound Name	Sample Quantitation Limit (ug/l)	Compound of Concern	Result (ug/l)	GWQC (ug/l)	Exceeds Criteria
Trip Blank	11/28/95	12/11/95	VOLATILE ORGANICS (Continued)					
			4-Chlorotoluene	0.50	--	ND	--	--
			1,3,5-Trimethylbenzene	0.50	--	ND	--	--
			tert-Butylbenzene	0.50	--	ND	--	--
			1,2,4-Trimethylbenzene	0.50	--	ND	--	--
			sec-Butylbenzene	0.50	--	ND	--	--
			1,3-Dichlorobenzene	0.50	--	ND	600	--
			4-Isopropyltoluene	0.50	--	ND	75	--
			1,4-Dichlorobenzene	0.50	--	ND	--	--
			1,2-Dichlorobenzene	0.50	--	ND	600	--
			N-Butylbenzene	0.50	--	ND	--	--
			1,2-Dibromo-3-chloropropane	0.50	--	ND	NA	--
			1,2,4-Trichlorobenzene	0.50	--	ND	9	--
			Hexachlorobutadiene	0.50	--	ND	1	--
			Naphthalene	0.50	--	ND	--	--
			1,2,3-Trichlorobenzene	0.50	--	ND	--	--
			Methy-tertiary butyl ether	0.50	--	ND	--	--
			tertiary-Butyl alcohol	2.0	--	ND	--	--
			VOLATILE TICS:					
			Furan, tetraydro-	--	--	3 J	--	--
			Column Bleed	--	--	1 J	--	--
			TOTAL TICS:	--	--	4	--	--

TABLE 3

GROUNDWATER SAMPLING RESULTS
BUILDING 290, MAIN POST, FIELD BLANK
FT. MONMOUTH, NEW JERSEY

Sample ID	Sample Date	Analysis Date	Compound Name	Sample Quantitation Limit (ug/l)	Compound of Concern	Result (ug/l)	GWQC (ug/l)	Exceeds Criteria
Field Blank	11/28/95	12/11/95	VOLATILE ORGANICS					
			Dichlorodifluoromethane	0.50	--	ND	--	--
			Chloromethane	0.50	--	ND	--	--
			Vinyl Chloride	0.50	--	ND	--	--
			Bromomethane	0.50	--	ND	5	--
			Chloroethane	0.50	--	ND	--	--
			Trichlorofluoromethane	0.50	--	ND	--	--
			1,1-Dichloroethene	0.50	--	ND	2	--
			Methylene Chloride	3.1	--	3.1 B	100	--
			1,2-Dichloroethene (trans)	0.50	--	ND	2	--
			1,1 Dichloroethane	0.50	--	ND	70	--
			2,2-Dichloropropane	0.50	--	ND	--	--
			cis-1,2-Dichloroethene	0.50	--	ND	--	--
			Bromochloromethane	0.50	--	ND	10	--
			Chloroform	0.50	--	ND	6	--
			1,1,1-Trichloroethane	0.50	--	ND	--	--
			Carbon Tetrachloride	0.50	--	ND	2	--
			1,1-Dichloropropene	0.50	--	ND	30	--
			Benzene	0.50	--	ND	--	--
			1,2-Dichloroethane	0.50	--	ND	2	--
			Trichloroethene	0.50	--	ND	1	--
			1,2-Dichloropropane	0.50	--	ND	1	--
			Dibromomethane	0.50	--	ND	NA	--
			Bromodichloromethane	0.50	--	ND	--	--
			cis-1,3-Dichloropropene	0.50	--	ND	1	--
			Toluene	0.50	--	ND	10	--
			trans-1,3-Dichloropropene	0.50	--	ND	3	--
			1,1,2-Trichloroethane	0.50	--	ND	1	--
			Tetrachloroethene	0.50	--	ND	NA	--
			1,3-Dichloropropane	0.50	--	ND	4	--
			Dibromochloromethane	0.50	--	ND	10	--
			1,2-Dibromomethane	0.50	--	ND	1	--
			Chlorobenzene	0.50	--	ND	2	--
			1,1,2,2-Tetrachloroethane	0.50	--	ND	1,000	--
			Ethylbenzene	0.50	--	ND	--	--
			Xylenes (Total)	0.50	--	ND	4	--
			Styrene	0.50	--	ND	700	--
			Bromoform	0.50	--	ND	40	--
			Isopropylbenzene	0.50	--	ND	100	--
			Bromobenzene	0.50	--	ND	--	--
			1,1,2,2-Tetrachloroethane	0.50	--	ND	--	--
			1,2,3-Trichloropropane	0.50	--	ND	--	--
			n-Propylbenzene	0.50	--	ND	--	--
			2-Chlorotoluene	0.50	--	ND	--	--

TABLE 3

GROUNDWATER SAMPLING RESULTS
BUILDING 290, MAIN POST, FIELD BLANK
FORT MONMOUTH, NEW JERSEY

Sample ID	Sample Date	Analysis Date	Compound Name	Sample Quantitation Limit (ug/l)	Compound of Concern	Result (ug/l)	GWQC (ug/l)	Exceeds Criteria
Field Blank	11/28/95	12/11/95	VOLATILE ORGANICS (Continued)					
			4-Chlorotoluene	0.50	--	ND	--	--
			1,3,5-Trimethylbenzene	0.50	--	ND	--	--
			tert-Butylbenzene	0.50	--	ND	--	--
			1,2,4-Trimethylbenzene	0.50	--	ND	--	--
			sec-Butylbenzene	0.50	--	ND	--	--
			1,3-Dichlorobenzene	0.50	--	ND	600	--
			4-Isopropyltoluene	0.50	--	ND	75	--
			1,4-Dichlorobenzene	0.50	--	ND	--	--
			1,2-Dichlorobenzene	0.50	--	ND	600	--
			N-Butylbenzene	0.50	--	ND	--	--
			1,2-Dibromo-3-chloropropane	0.50	--	ND	NA	--
			1,2,4-Trichlorobenzene	0.50	--	ND	9	--
			Hexachlorobutadiene	0.50	--	ND	1	--
			Naphthalene	0.50	--	ND	--	--
			1,2,3-Trichlorobenzene	0.50	--	ND	--	--
			Methy-tertiary butyl ether	0.50	--	ND	--	--
			tertiary-Butyl alcohol	2.0	--	ND	--	--
			VOLATILE TICS:					
			Colume Bleed	--	--	1 J	--	--
			Unknown	--	--	3 J	--	--
			Colume Bleed	--	--	1 J	--	--
			Unknown Hydrocarbon	--	--	1 J	--	--
			TOTAL TICS:	--	--	6	--	--

TABLE 3

GROUNDWATER SAMPLING RESULTS
BUILDING 290, MAIN POST, FIELD BLANK
FORT MONMOUTH, NEW JERSEY

Sample ID	Sample Date	Analysis Date	Compound Name	Sample Quantitation Limit (ug/l)	Compound of Concern	Result (ug/l)	GWQC (ug/l)	Exceeds Criteria
FIELD BLANK	11/28/95	12/11/95	SEMIVOLATILES					
			Phenol	10	--	ND	4000	--
			bis(2 chloroethyl)Ether	10	--	ND	20	--
			2-Chlorophenol	10	--	ND	40	--
			1,3-Dichlorobenzene	10	--	ND	600	--
			1,4-Dichlorobenzene	10	--	ND	75	--
			1,2-Dichlorobenzene	10	--	ND	600	--
			2-Methylphenol	10	--	ND	--	--
			bis(2-chloroisopropyl)Ether	10	--	ND	300	--
			4-Methylphenol	10	-	ND	--	--
			N-Nitroso-Di-n-propylamine	10	--	ND	20	--
			Hexachloroethane	10	--	ND	10	--
			Nitrobenzene	10	--	ND	10	--
			Isophorone	10	--	ND	100	--
			2-Nitrophenol	10	--	ND	--	--
			2,4-Dimethylphenol	10	--	ND	100	--
			bis(2-Chloroethoxy)Methane	10	--	ND	--	--
			2,4-Dichlorophenol	10	--	ND	20	--
			1,2,4-Trichlorobenzene	10	--	ND	9	--
			Naphthalene	10	--	ND	--	--
			4-Chloroaniline	10	--	ND	--	--
			Hexachlorobutadiene	10	--	ND	1	--
			4-Chloro-3-methylphenol	10	--	ND	--	--
			2-methylnaphthanene	10	--	ND	--	--
			Hexachlorocyclopentadiene	10	--	ND	50	--
			2,4,6-Trichlorophenol	10	--	ND	20	--
			2,4,5-Trichlorophenol	25	--	ND	700	--
			2-Chloronaphthalene	10	--	ND	--	--
			2-Nitroanline	25	--	ND	--	--
			Dimethyl Phthalate	10	--	ND	--	--
			Acenaphthylene	10	--	ND	NA	--
			2,6-Dinitrotoluene	10	--	ND	NA	--
			3-Nitroaniline	25	--	ND	--	--
			Acenaphthene	10	--	ND	400	--
			2,4-Dinitrophenol	25	--	ND	40	--
			4-Nitrophenol	25	--	ND	--	--
			Dibenzofuran	10	--	ND	--	--
			2,4-Dinitrotoluene	10	--	ND	10	--
			Diethylphthalate	10	--	ND	5,000	--
			Fluorene	10	--	ND	300	--
			4-Chlorophenyl-phenlyether	10	--	ND	--	--
			4-Nitroaniline	25	--	ND	--	--
			4,6-Dinitro-2-methylphenol	25	--	ND	--	--
			N-Nitrosodiphenylamine	10	--	ND	20	--

TABLE 3

GROUNDWATER SAMPLING RESULTS
BUILDING 290, MAIN POST, FIELD BLANK
FORT MONMOUTH, NEW JERSEY

Sample ID	Sample Date	Analysis Date	Compound Name	Sample Quantitation Limit (ug/l)	Compound of Concern	Result (ug/l)	GWQC (ug/l)	Exceeds Criteria
FIELD BLANK	11/28/95	12/11/95	SEMIVOLATILES (Continued)					
			4-Bromophenyl-phenylether	10	--	ND	--	--
			Hexachlorobenzene	10	--	ND	10	--
			Pentachlorophenol	25	--	ND	1	--
			Phenanthrene	10	--	ND	NA	--
			Anthracene	10	--	ND	2,000	--
			Carbazole	10	--	ND	--	--
			Di-n-butylphthalate	10	--	ND	900	--
			Fluoranthene	10	--	ND	300	--
			Pyrene	10	--	ND	200	--
			Butylbenzylphthalate	10	--	ND	100	--
			Benzo(a)Anthracene	10	--	ND	NA	--
			3,3-Dichlorobenzidine	20	--	ND	60	--
			Chrysene	10	--	ND	NA	--
			bis(2-Ethylhexyl)Phthalate	10	--	ND	30	--
			Di-n-Octyl Phthalate	10	--	ND	100	--
			Benzo(b)Fluoranthene	10	--	ND	NA	--
			Benzo(k)Fluoranthene	10	--	ND	NA	--
			Benzo(a)Pyrene	10	--	ND	NA	--
			Indeno(1,2,3-cd)pyrene	10	--	ND	NA	--
			Dibenzo(a,h)anthracene	10	--	ND	NA	--
			Benzo(g,h,i)perylene	10	--	ND	NA	--
			SEMIVOLATILE TICS:					
			Benzene,1-chloro-2-methyl-	--	--	9 J	--	--
			Benzene,1-chloro-4-methyl-	--	--	2 J	--	--
			Unknown Hydrocarbon		--	5 J	--	--
			Hexadecanoic		--	12 J	--	--
			Unknown Hydrocarbon		--	2 J	--	--
			TOTAL TICS:	--	--	35	--	--

TABLE 3

GROUNDWATER SAMPLING RESULTS
BUILDING 290, MAIN POST, FIELD BLANK
FORT MONMOUTH, NEW JERSEY

Sample ID	Sample Date	Analysis Date	Compound Name	Sample Quantitation Limit (ug/l)	Compound of Concern	Result (ug/l)	GWQC (ug/l)	Exceeds Criteria
Field Blank	11/28/95	12/07/95	METALS					
			Silver	110	--	110	NA	--
			Aluminum	200	--	ND	200	--
			Arsenic	5	--	ND	8	--
			Barium	20	--	ND	2,000	--
			Beryllium	5	--	ND	20	--
			Calcium	400	--	ND	--	--
			Cadmium	10	--	ND	4	--
			Cobalt	50	--	ND	--	--
			Chromium	50	--	ND	100	--
			Copper	50	--	ND	1,000	--
			Iron	100	--	ND	300	--
			Mercury	1	--	ND	2	--
			Potassium	3,000	--	ND	--	--
			Magnesium	200	--	ND	--	--
			Manganese	33	--	32	50	--
			Sodium	400	--	ND	50,000	--
			Nickel	50	--	ND	100	--
			Lead	2.8	--	2.8	10	--
			Antimony	5	--	ND	20	--
			Selenium	5	--	ND	50	--
			Thallium	2	--	ND	10	--
			Vanadium	50	--	ND	--	--
			Zinc	27	--	27	5,000	--
			Cyanide (Total)	10	--	ND	200	--

TABLE 3

GROUNDWATER SAMPLING RESULTS
BUILDING 290, MAIN POST, FIELD BLANK
FORT MONMOUTH, NEW JERSEY

Sample ID	Sample Date	Analysis Date	Compound Name	Sample Quantitation Limit (ug/l)	Compound of Concern	Result (ug/l)	GWQC (ug/l)	Exceeds Criteria
Field Blank	11/28/95	12/07/95	PESTICIDES/PCBs					
			alpha-BHC	0.02	--	ND	0.02	--
			beta-BHC	0.04	--	ND	0.2	--
			delta-BHC	0.02	--	ND	--	--
			gamma-BHC (Lindane)	0.03	--	ND	0.2	--
			Heptachlor	0.02	--	ND	0.4	--
			Aldrin	0.04	--	ND	0.04	--
			Heptachlor epoxide	0.05	--	ND	0.2	--
			Endosulfan I	0.04	--	ND	0.4	--
			Dieldrin	0.03	--	ND	0.03	--
			4,4'-DDE	0.04	--	ND	0.1	--
			Endrin	0.04	--	ND	2.0	--
			Endosulfan II	0.04	--	ND	0.4	--
			4,4'-DDD	0.04	--	ND	0.1	--
			Endosulfan sulfate	0.08	--	ND	0.4	--
			4,4'-DDT	0.04	--	ND	0.1	--
			Endrin aldehyde	0.1	--	ND	--	--
			Chlordane	0.1	--	ND	0.5	--
			Toxaphene	1.0	--	ND	3.0	--
			Aroclor-1016	1.0	--	ND	--	--
			Aroclor-1221	1.0	--	ND	--	--
			Aroclor-1232	1.0	--	ND	--	--
			Aroclor-1242	1.0	--	ND	--	--
			Aroclor-1248	1.0	--	ND	--	--
			Aroclor-1254	1.0	--	ND	--	--
			Aroclor-1260	1.0	--	ND	--	--

TABLE 3

GROUNDWATER SAMPLING RESULTS
BUILDING 290, MAIN POST, MW-1
FT. MONMOUTH, NEW JERSEY

Sample ID	Sample Date	Analysis Date	Compound Name	Sample Quantitation Limit (ug/l)	Compound of Concern	Result (ug/l)	GWQC (ug/l)	Exceeds Criteria
MW-1	12/18/95	12/29/95	VOLATILE ORGANICS					
			Dichlorodifluoromethane	0.50	--	ND	--	--
			Chloromethane	0.50	--	ND	--	--
			Vinyl Chloride	0.50	--	ND	--	--
			Bromomethane	0.50	--	ND	5	--
			Chloroethane	0.50	--	ND	--	--
			Trichlorofluoromethane	0.50	--	ND	--	--
			1,1-Dichloroethene	0.50	--	ND	2	--
			Methylene Chloride	1.1	--	1.1 B	100	--
			1,2-Dichloroethene (trans)	0.50	--	ND	2	--
			1,1 Dichloroethane	0.50	--	ND	70	--
			2,2-Dichloropropane	0.50	--	ND	--	--
			cis-1,2-Dichloroethene	1.1	--	1.1	--	--
			Bromochloromethane	0.50	--	ND	10	--
			Chloroform	0.50	--	ND	6	--
			1,1,1-Trichloroethane	0.50	--	ND	--	--
			Carbon Tetrachloride	0.50	--	ND	2	--
			1,1-Dichloropropene	0.50	--	ND	30	--
			Benzene	0.50	--	ND	--	--
			1,2-Dichloroethane	0.50	--	ND	2	--
			Trichloroethene	0.50	--	ND	1	--
			1,2-Dichloropropane	0.50	--	ND	1	--
			Dibromomethane	0.50	--	ND	NA	--
			Bromodichloromethane	0.50	--	ND	--	--
			cis-1,3-Dichloropropene	0.50	--	ND	1	--
			Toluene	0.50	--	ND	10	--
			trans-1,3-Dichloropropene	0.50	--	ND	3	--
			1,1,2-Trichloroethane	0.50	--	ND	1	--
			Tetrachloroethene	0.50	--	ND	NA	--
			1,3-Dichloropropane	0.50	--	ND	4	--
			Dibromochloromethane	0.50	--	ND	10	--
			1,2-Dibromomethane	0.50	--	ND	1	--
			Chlorobenzene	0.50	--	ND	2	--
			1,1,2,2-Tetrachloroethane	0.50	--	ND	1,000	--
			Ethylbenzene	0.50	--	ND	--	--
			Xylenes (Total)	0.50	--	ND	4	--
			Styrene	0.50	--	ND	700	--
			Bromoform	0.50	--	ND	40	--
			Isopropylbenzene	0.50	--	ND	100	--
			Bromobenzene	0.50	--	ND	--	--
			1,1,2,2-Tetrachloroethane	0.50	--	ND	--	--
			1,2,3-Trichloropropane	0.50	--	ND	--	--
			n-Propylbenzene	0.50	--	ND	--	--
			2-Chlorotoluene	0.50	--	ND	--	--

TABLE 3

GROUNDWATER SAMPLING RESULTS
BUILDING 290, MAIN POST, MW-1
FORT MONMOUTH, NEW JERSEY

Sample ID	Sample Date	Analysis Date	Compound Name	Sample Quantitation Limit (ug/l)	Compound of Concern	Result (ug/l)	GWQC (ug/l)	Exceeds Criteria
MW-1	12/18/95	12/29/95	VOLATILE ORGANICS (Continued)					
			4-Chlorotoluene	0.50	--	ND	--	--
			1,3,5-Trimethylbenzene	0.50	--	ND	--	--
			tert-Butylbenzene	0.50	--	ND	--	--
			1,2,4-Trimethylbenzene	0.50	--	ND	--	--
			sec-Butylbenzene	0.50	--	ND	--	--
			1,3-Dichlorobenzene	0.50	--	ND	600	--
			4-Isopropyltoluene	0.50	--	ND	75	--
			1,4-Dichlorobenzene	0.50	--	ND	--	--
			1,2-Dichlorobenzene	0.50	--	ND	600	--
			N-Butylbenzene	0.50	--	ND	--	--
			1,2-Dibromo-3-chloropropane	0.50	--	ND	NA	--
			1,2,4-Trichlorobenzene	0.50	--	ND	9	--
			Hexachlorobutadiene	0.50	--	ND	1	--
			Naphthalene	0.50	--	ND	--	--
			1,2,3-Trichlorobenzene	0.50	--	ND	--	--
			Methy-tertiary butyl ether	0.50	--	ND	--	--
			tertiary-Butyl alcohol	2.0	--	ND	--	--
			VOLATILE TICS:					
			Column Bleed	--	--	3 J	--	--
			Column Bleed	--	--	5 J	--	--
			TOTAL TICS:	--	--	8	--	--

TABLE 3

GROUNDWATER SAMPLING RESULTS
BUILDING 290, MAIN POST, MW-1
FORT MONMOUTH, NEW JERSEY

Sample ID	Sample Date	Analysis Date	Compound Name	Sample Quantitation Limit (ug/l)	Compound of Concern	Result (ug/l)	GWQC (ug/l)	Exceeds Criteria
MW-1	12/18/95	12/26/95	SEMIVOLATILES					
			Phenol	10	--	ND	4000	--
			bis(2 chloroethyl)Ether	10	--	ND	20	--
			2-Chlorophenol	10	--	ND	40	--
			1,3-Dichlorobenzene	10	--	ND	600	--
			1,4-Dichlorobenzene	10	--	ND	75	--
			1,2-Dichlorobenzene	10	--	ND	600	--
			2-Methylphenol	10	--	ND	--	--
			bis(2-chloroisopropyl)Ether	10	--	ND	300	--
			4-Methylphenol	10	-	ND	--	--
			N-Nitroso-Di-n-propylamine	10	--	ND	20	--
			Hexachloroethane	10	--	ND	10	--
			Nitrobenzene	10	--	ND	10	--
			Isophorone	10	--	ND	100	--
			2-Nitrophenol	10	--	ND	--	--
			2,4-Dimethylphenol	10	--	ND	100	--
			bis(2-Chloroethoxy)Methane	10	--	ND	--	--
			2,4-Dichlorophenol	10	--	ND	20	--
			1,2,4-Trichlorobenzene	10	--	ND	9	--
			Naphthalene	10	--	ND	--	--
			4-Chloroaniline	10	--	ND	--	--
			Hexachlorobutadiene	10	--	ND	1	--
			4-Chloro-3-methylphenol	10	--	ND	--	--
			2-methylnaphthanene	10	--	ND	--	--
			Hexachlorocyclopentadiene	10	--	ND	50	--
			2,4,6-Trichlorophenol	10	--	ND	20	--
			2,4,5-Trichlorophenol	25	--	ND	700	--
			2-Chloronaphthalene	10	--	ND	--	--
			2-Nitroanline	25	--	ND	--	--
			Dimethyl Phthalate	10	--	ND	--	--
			Acenaphthylene	10	--	ND	NA	--
			2,6-Dinitrotoluene	10	--	ND	NA	--
			3-Nitroaniline	25	--	ND	--	--
			Acenaphthene	10	--	ND	400	--
			2,4-Dinitrophenol	25	--	ND	40	--
			4-Nitrophenol	25	--	ND	--	--
			Dibenzofuran	10	--	ND	--	--
			2,4-Dinitrotoluene	10	--	ND	10	--
			Diethylphthalate	10	--	ND	5,000	--
			Fluorene	10	--	ND	300	--
			4-Chlorophenyl-phenlyether	10	--	ND	--	--
			4-Nitroaniline	25	--	ND	--	--
			4,6-Dinitro-2-methylphenol	25	--	ND	--	--
			N-Nitrosodiphenylamine	10	--	ND	20	--

TABLE 3

GROUNDWATER SAMPLING RESULTS
BUILDING 290, MAIN POST, MW-1
FORT MONMOUTH, NEW JERSEY

Sample ID	Sample Date	Analysis Date	Compound Name	Sample Quantitation Limit (ug/l)	Compound of Concern	Result (ug/l)	GWQC (ug/l)	Exceeds Criteria
MW-1	12/18/95	12/26/95	SEMIVOLATILES (Continued)					
			4-Bromophenyl-phenylether	10	--	ND	--	--
			Hexachlorobenzene	10	--	ND	10	--
			Pentachlorophenol	25	--	ND	1	--
			Phenanthrene	10	--	ND	NA	--
			Anthracene	10	--	ND	2,000	--
			Carbazole	10	--	ND	--	--
			Di-n-butylphthalate	10	--	ND	900	--
			Fluoranthene	10	--	ND	300	--
			Pyrene	10	--	ND	200	--
			Butylbenzylphthalate	10	--	ND	100	--
			Benzo(a)Anthracene	10	--	ND	NA	--
			3,3-Dichlorobenzidine	20	--	ND	60	--
			Chrysene	10	--	ND	NA	--
			bis(2-Ethylhexyl)Phthalate	10	--	ND	30	--
			Di-n-Octyl Phthalate	10	--	ND	100	--
			Benzo(b)Fluoranthene	10	--	ND	NA	--
			Benzo(k)Fluoranthene	10	--	ND	NA	--
			Benzo(a)Pyrene	10	--	ND	NA	--
			Indeno(1,2,3-cd)pyrene	10	--	ND	NA	--
			Dibenzo(a,h)anthracene	10	--	ND	NA	--
			Benzo(g,h,i)perylene	10	--	ND	NA	--
			SEMIVOLATILE TICS:					
			NONE FOUND	--	--	--	--	--

TABLE 3

GROUNDWATER SAMPLING RESULTS
BUILDING 290, MAIN POST, MW-1
FORT MONMOUTH, NEW JERSEY

Sample ID	Sample Date	Analysis Date	Compound Name	Sample Quantitation Limit (ug/l)	Compound of Concern	Result (ug/l)	GWQC (ug/l)	Exceeds Criteria
MW-1	12/18/95	12/29/95	METALS					
			Silver	50	--	ND	NA	--
			Aluminum	2,200	--	2,200	200	YES
			Arsenic	1	--	ND	8	--
			Barium	55	--	55	2,000	--
			Beryllium	5	--	ND	20	--
			Calcium	4,100	--	4,100	--	--
			Cadmium	10	--	ND	4	--
			Cobalt	50	--	ND	--	--
			Chromium	50	--	ND	100	--
			Copper	60	--	60	1,000	--
			Iron	6,400	--	6,400	300	YES
			Mercury	0.2	--	ND	2	--
			Potassium	5,000	--	5,000	--	--
			Magnesium	4,500	--	4,500	--	--
			Manganese	23	--	23	50	--
			Sodium	9,200	--	9,200	50,000	--
			Nickel	50	--	ND	100	--
			Lead	2	--	ND	10	--
			Antimony	5	--	ND	20	--
			Selenium	5	--	ND	50	--
			Thallium	2	--	ND	10	--
			Vanadium	50	--	ND	--	--
			Zinc	180	--	180	5,000	--
			Cyanide (Total)	10	--	ND	200	--

TABLE 3

GROUNDWATER SAMPLING RESULTS
BUILDING 290, MAIN POST, MW-1
FORT MONMOUTH, NEW JERSEY

Sample ID	Sample Date	Analysis Date	Compound Name	Sample Quantitation Limit (ug/l)	Compound of Concern	Result (ug/l)	GWQC (ug/l)	Exceeds Criteria
MW-1	12/18/95	12/29/95	PESTICIDES/PCBs					
			alpha-BHC	0.02	--	ND	0.02	--
			beta-BHC	0.04	--	ND	0.2	--
			delta-BHC	0.02	--	ND	--	--
			gamma-BHC (Lindane)	0.03	--	ND	0.2	--
			Heptachlor	0.02	--	ND	0.4	--
			Aldrin	0.04	--	ND	0.04	--
			Heptachlor epoxide	0.05	--	ND	0.2	--
			Endosulfan I	0.04	--	ND	0.4	--
			Dieldrin	0.03	--	ND	0.03	--
			4,4'-DDE	0.04	--	ND	0.1	--
			Endrin	0.04	--	ND	2.0	--
			Endosulfan II	0.04	--	ND	0.4	--
			4,4'-DDD	0.04	--	ND	0.1	--
			Endosulfan sulfate	0.08	--	ND	0.4	--
			4,4'-DDT	0.04	--	ND	0.1	--
			Endrin aldehyde	0.1	--	ND	--	--
			Chlordane	0.1	--	ND	0.5	--
			Toxaphene	1.0	--	ND	3.0	--
			Aroclor-1016	1.0	--	ND	--	--
			Aroclor-1221	1.0	--	ND	--	--
			Aroclor-1232	1.0	--	ND	--	--
			Aroclor-1242	1.0	--	ND	--	--
			Aroclor-1248	1.0	--	ND	--	--
			Aroclor-1254	1.0	--	ND	--	--
			Aroclor-1260	1.0	--	ND	--	--

TABLE 3

GROUNDWATER SAMPLING RESULTS
BUILDING 290, MAIN POST, TRIP BLANK
FT. MONMOUTH, NEW JERSEY

Sample ID	Sample Date	Analysis Date	Compound Name	Sample Quantitation Limit (ug/l)	Compound of Concern	Result (ug/l)	GWQC (ug/l)	Exceeds Criteria
Trip Blank	12/18/95	12/29/95	VOLATILE ORGANICS					
			Dichlorodifluoromethane	0.50	--	ND	--	--
			Chloromethane	0.50	--	ND	--	--
			Vinyl Chloride	0.50	--	ND	--	--
			Bromomethane	0.50	--	ND	5	--
			Chloroethane	0.50	--	ND	--	--
			Trichlorofluoromethane	0.50	--	ND	--	--
			1,1-Dichloroethene	0.50	--	ND	2	--
			Methylene Chloride	1.4	--	1.4 B	100	--
			1,2-Dichloroethene (trans)	0.50	--	ND	2	--
			1,1 Dichloroethane	0.50	--	ND	70	--
			2,2-Dichloropropane	0.50	--	ND	--	--
			cis-1,2-Dichloroethene	0.50	--	ND	--	--
			Bromochloromethane	0.50	--	ND	10	--
			Chloroform	0.50	--	ND	6	--
			1,1,1-Trichloroethane	0.50	--	ND	--	--
			Carbon Tetrachloride	0.50	--	ND	2	--
			1,1-Dichloropropene	0.50	--	ND	30	--
			Benzene	0.50	--	ND	--	--
			1,2-Dichloroethane	0.50	--	ND	2	--
			Trichloroethene	0.50	--	ND	1	--
			1,2-Dichloropropane	0.50	--	ND	1	--
			Dibromomethane	0.50	--	ND	NA	--
			Bromodichloromethane	0.50	--	ND	--	--
			cis-1,3-Dichloropropene	0.50	--	ND	1	--
			Toluene	0.50	--	ND	10	--
			trans-1,3-Dichloropropene	0.50	--	ND	3	--
			1,1,2-Trichloroethane	0.50	--	ND	1	--
			Tetrachloroethene	0.50	--	ND	NA	--
			1,3-Dichloropropane	0.50	--	ND	4	--
			Dibromochloromethane	0.50	--	ND	10	--
			1,2-Dibromomethane	0.50	--	ND	1	--
			Chlorobenzene	0.50	--	ND	2	--
			1,1,2,2-Tetrachloroethane	0.50	--	ND	1,000	--
			Ethylbenzene	0.50	--	ND	--	--
			Xylenes (Total)	0.50	--	ND	4	--
			Styrene	0.50	--	ND	700	--
			Bromoform	0.50	--	ND	40	--
			Isopropylbenzene	0.50	--	ND	100	--
			Bromobenzene	0.50	--	ND	--	--
			1,1,2,2-Tetrachloroethane	0.50	--	ND	--	--
			1,2,3-Trichloropropane	0.50	--	ND	--	--
			n-Propylbenzene	0.50	--	ND	--	--
			2-Chlorotoluene	0.50	--	ND	--	--

TABLE 3

GROUNDWATER SAMPLING RESULTS
BUILDING 290, MAIN POST, TRIP BLANK
FORT MONMOUTH, NEW JERSEY

Sample ID	Sample Date	Analysis Date	Compound Name	Sample Quantitation Limit (ug/l)	Compound of Concern	Result (ug/l)	GWQC (ug/l)	Exceeds Criteria
Trip Blank	12/18/95	12/29/95	VOLATILE ORGANICS (Continued)					
			4-Chlorotoluene	0.50	--	ND	--	--
			1,3,5-Trimethylbenzene	0.50	--	ND	--	--
			tert-Butylbenzene	0.50	--	ND	--	--
			1,2,4-Trimethylbenzene	0.50	--	ND	--	--
			sec-Butylbenzene	0.50	--	ND	--	--
			1,3-Dichlorobenzene	0.50	--	ND	600	--
			4-Isopropyltoluene	0.50	--	ND	75	--
			1,4-Dichlorobenzene	0.50	--	ND	--	--
			1,2-Dichlorobenzene	0.50	--	ND	600	--
			N-Butylbenzene	0.50	--	ND	--	--
			1,2-Dibromo-3-chloropropane	0.50	--	ND	NA	--
			1,2,4-Trichlorobenzene	0.50	--	ND	9	--
			Hexachlorobutadiene	0.50	--	ND	1	--
			Naphthalene	0.50	--	ND	--	--
			1,2,3-Trichlorobenzene	0.50	--	ND	--	--
			Methy-tertiary butyl ether	0.50	--	ND	--	--
			tertiary-Butyl alcohol	2.0	--	ND	--	--
			VOLATILE TICS:					
			Furan, tetraydro-	--	--	2 J	--	--
			Column Bleed	--	--	2 J	--	--
			Column Bleed	--	--	1 J	--	--

TABLE 3

GROUNDWATER SAMPLING RESULTS
BUILDING 290, MAIN POST, FIELD BLANK
FT. MONMOUTH, NEW JERSEY

Sample ID	Sample Date	Analysis Date	Compound Name	Sample Quantitation Limit (ug/l)	Compound of Concern	Result (ug/l)	GWQC (ug/l)	Exceeds Criteria
Field Blank	12/18/95	12/29/95	VOLATILE ORGANICS					
			Dichlorodifluoromethane	0.50	--	ND	--	--
			Chloromethane	0.50	--	ND	--	--
			Vinyl Chloride	0.50	--	ND	--	--
			Bromomethane	0.50	--	ND	5	--
			Chloroethane	0.50	--	ND	--	--
			Trichlorofluoromethane	0.50	--	ND	--	--
			1,1-Dichloroethene	0.50	--	ND	2	--
			Methylene Chloride	1.4	--	1.4 B	100	--
			1,2-Dichloroethene (trans)	0.50	--	ND	2	--
			1,1 Dichloroethane	0.50	--	ND	70	--
			2,2-Dichloropropane	0.50	--	ND	--	--
			cis-1,2-Dichloroethene	0.50	--	ND	--	--
			Bromochloromethane	0.50	--	ND	10	--
			Chloroform	0.50	--	ND	6	--
			1,1,1-Trichloroethane	0.50	--	ND	--	--
			Carbon Tetrachloride	0.50	--	ND	2	--
			1,1-Dichloropropene	0.50	--	ND	30	--
			Benzene	0.50	--	ND	--	--
			1,2-Dichloroethane	0.50	--	ND	2	--
			Trichloroethene	0.50	--	ND	1	--
			1,2-Dichloropropane	0.50	--	ND	1	--
			Dibromomethane	0.50	--	ND	NA	--
			Bromodichloromethane	0.50	--	ND	--	--
			cis-1,3-Dichloropropene	0.50	--	ND	1	--
			Toluene	0.50	--	ND	10	--
			trans-1,3-Dichloropropene	0.50	--	ND	3	--
			1,1,2-Trichloroethane	0.50	--	ND	1	--
			Tetrachloroethene	0.50	--	ND	NA	--
			1,3-Dichloropropane	0.50	--	ND	4	--
			Dibromochloromethane	0.50	--	ND	10	--
			1,2-Dibromomethane	0.50	--	ND	1	--
			Chlorobenzene	0.50	--	ND	2	--
			1,1,2,2-Tetrachloroethane	0.50	--	ND	1,000	--
			Ethylbenzene	0.50	--	ND	--	--
			Xylenes (Total)	0.50	--	ND	4	--
			Styrene	0.50	--	ND	700	--
			Bromoform	0.50	--	ND	40	--
			Isopropylbenzene	0.50	--	ND	100	--
			Bromobenzene	0.50	--	ND	--	--
			1,1,2,2-Tetrachloroethane	0.50	--	ND	--	--
			1,2,3-Trichloropropane	0.50	--	ND	--	--
			n-Propylbenzene	0.50	--	ND	--	--
			2-Chlorotoluene	0.50	--	ND	--	--

TABLE 3

GROUNDWATER SAMPLING RESULTS
BUILDING 290, MAIN POST, FIELD BLANK
FORT MONMOUTH, NEW JERSEY

Sample ID	Sample Date	Analysis Date	Compound Name	Sample Quantitation Limit (ug/l)	Compound of Concern	Result (ug/l)	GWQC (ug/l)	Exceeds Criteria
Field Blank	12/18/95	12/29/95	VOLATILE ORGANICS (Continued)					
			4-Chlorotoluene	0.50	--	ND	--	--
			1,3,5-Trimethylbenzene	0.50	--	ND	--	--
			tert-Butylbenzene	0.50	--	ND	--	--
			1,2,4-Trimethylbenzene	0.50	--	ND	--	--
			sec-Butylbenzene	0.50	--	ND	--	--
			1,3-Dichlorobenzene	0.50	--	ND	600	--
			4-Isopropyltoluene	0.50	--	ND	75	--
			1,4-Dichlorobenzene	0.50	--	ND	--	--
			1,2-Dichlorobenzene	0.50	--	ND	600	--
			N-Butylbenzene	0.50	--	ND	--	--
			1,2-Dibromo-3-chloropropane	0.50	--	ND	NA	--
			1,2,4-Trichlorobenzene	0.50	--	ND	9	--
			Hexachlorobutadiene	0.50	--	ND	1	--
			Naphthalene	0.50	--	ND	--	--
			1,2,3-Trichlorobenzene	0.50	--	ND	--	--
			Methy-tertiary butyl ether	0.50	--	ND	--	--
			tertiary-Butyl alcohol	2.0	--	ND	--	--
			VOLATILE TICS:					
			Column Bleed	--	--	1 J	--	--
			Furan, tetrahydro-	--	--	2 J	--	--

TABLE 3

GROUNDWATER SAMPLING RESULTS
BUILDING 290, MAIN POST, FIELD BLANK
FORT MONMOUTH, NEW JERSEY

Sample ID	Sample Date	Analysis Date	Compound Name	Sample Quantitation Limit (ug/l)	Compound of Concern	Result (ug/l)	GWQC (ug/l)	Exceeds Criteria
Field Blank	12/18/95	12/26/95	SEMIVOLATILES					
			Phenol	10	--	ND	4000	--
			bis(2 chloroethyl)Ether	10	--	ND	20	--
			2-Chlorophenol	10	--	ND	40	--
			1,3-Dichlorobenzene	10	--	ND	600	--
			1,4-Dichlorobenzene	10	--	ND	75	--
			1,2-Dichlorobenzene	10	--	ND	600	--
			2-Methylphenol	10	--	ND	--	--
			bis(2-chloroisopropyl)Ether	10	--	ND	300	--
			4-Methylphenol	10	-	ND	--	--
			N-Nitroso-Di-n-propylamine	10	--	ND	20	--
			Hexachloroethane	10	--	ND	10	--
			Nitrobenzene	10	--	ND	10	--
			Isophorone	10	--	ND	100	--
			2-Nitrophenol	10	--	ND	--	--
			2,4-Dimethylphenol	10	--	ND	100	--
			bis(2-Chloroethoxy)Methane	10	--	ND	--	--
			2,4-Dichlorophenol	10	--	ND	20	--
			1,2,4-Trichlorobenzene	10	--	ND	9	--
			Naphthalene	10	--	ND	--	--
			4-Chloroaniline	10	--	ND	--	--
			Hexachlorobutadiene	10	--	ND	1	--
			4-Chloro-3-methylphenol	10	--	ND	--	--
			2-methylnaphthanene	10	--	ND	--	--
			Hexachlorocyclopentadiene	10	--	ND	50	--
			2,4,6-Trichlorophenol	10	--	ND	20	--
			2,4,5-Trichlorophenol	25	--	ND	700	--
			2-Chloronaphthalene	10	--	ND	--	--
			2-Nitroanline	25	--	ND	--	--
			Dimethyl Phthalate	10	--	ND	--	--
			Acenaphthylene	10	--	ND	NA	--
			2,6-Dinitrotoluene	10	--	ND	NA	--
			3-Nitroaniline	25	--	ND	--	--
			Acenaphthene	10	--	ND	400	--
			2,4-Dinitrophenol	25	--	ND	40	--
			4-Nitrophenol	25	--	ND	--	--
			Dibenzofuran	10	--	ND	--	--
			2,4-Dinitrotoluene	10	--	ND	10	--
			Diethylphthalate	10	--	ND	5,000	--
			Fluorene	10	--	ND	300	--
			4-Chlorophenyl-phenlyether	10	--	ND	--	--
			4-Nitroaniline	25	--	ND	--	--
			4,6-Dinitro-2-methylphenol	25	--	ND	--	--
			N-Nitrosodiphenylamine	10	--	ND	20	--

TABLE 3

GROUNDWATER SAMPLING RESULTS
BUILDING 290, MAIN POST, FIELD BLANK
FORT MONMOUTH, NEW JERSEY

Sample ID	Sample Date	Analysis Date	Compound Name	Sample Quantitation Limit (ug/l)	Compound of Concern	Result (ug/l)	GWQC (ug/l)	Exceeds Criteria
Field Blank	12/18/95	12/26/95	SEMIVOLATILES (Continued)					
			4-Bromophenyl-phenylether	10	--	ND	--	--
			Hexachlorobenzene	10	--	ND	10	--
			Pentachlorophenol	25	--	ND	1	--
			Phenanthrene	10	--	ND	NA	--
			Anthracene	10	--	ND	2,000	--
			Carbazole	10	--	ND	--	--
			Di-n-butylphthalate	10	--	ND	900	--
			Fluoranthene	10	--	ND	300	--
			Pyrene	10	--	ND	200	--
			Butylbenzylphthalate	10	--	ND	100	--
			Benzo(a)Anthracene	10	--	ND	NA	--
			3,3-Dichlorobenzidine	20	--	ND	60	--
			Chrysene	10	--	ND	NA	--
			bis(2-Ethylhexyl)Phthalate	10	--	ND	30	--
			Di-n-Octyl Phthalate	10	--	ND	100	--
			Benzo(b)Fluoranthene	10	--	ND	NA	--
			Benzo(k)Fluoranthene	10	--	ND	NA	--
			Benzo(a)Pyrene	10	--	ND	NA	--
			Indeno(1,2,3-cd)pyrene	10	--	ND	NA	--
			Dibenzo(a,h)anthracene	10	--	ND	NA	--
			Benzo(g,h,i)perylene	10	--	ND	NA	--
			SEMIVOLATILE TICS:					
			Unkown Hydrocarbon	--	--	6 J	--	--

TABLE 3

GROUNDWATER SAMPLING RESULTS
BUILDING 290, MAIN POST, FIELD BLANK
FORT MONMOUTH, NEW JERSEY

Sample ID	Sample Date	Analysis Date	Compound Nane	Sample Quantitatio n Limit (ug/l)	Compound of Concern	Result (ug/l)	GWQC (ug/l)	Exceeds Criteria
Field Blank	12/18/95	12/29/95	METALS					
			Silver	50	--	ND	NA	--
			Aluminum	200	--	ND	200	--
			Arsenic	1	--	ND	8	--
			Barium	20	--	ND	2,000	--
			Beryllium	5	--	ND	20	--
			Calcium	400	--	ND	--	--
			Cadmium	10	--	ND	4	--
			Cobalt	50	--	ND	--	--
			Chromium	50	--	ND	100	--
			Copper	50	--	ND	1,000	--
			Iron	100	--	ND	300	--
			Mercury	0.2	--	ND	2	--
			Potassium	3,000	--	ND	--	--
			Magnesium	200	--	ND	--	--
			Manganese	20	--	ND	50	--
			Sodium	400	--	ND	50,000	--
			Nickel	50	--	ND	100	--
			Lead	2	--	ND	10	--
			Antimony	10	--	ND	20	--
			Selenium	5	--	ND	50	--
			Thallium	2	--	ND	10	--
			Vanadium	50	--	ND	--	--
			Zinc	65	--	65	5,000	--
			Cyanide (Total)	10	--	ND	200	--

TABLE 3

GROUNDWATER SAMPLING RESULTS
BUILDING 290, MAIN POST, FIELD BLANK
FORT MONMOUTH, NEW JERSEY

Sample ID	Sample Date	Analysis Date	Compound Nane	Sample Quantitation Limit (ug/l)	Compound of Concern	Result (ug/l)	GWQC (ug/l)	Exceeds Criteria
Field Blank	12/28/95	12/29/95	PESTICIDES/PCBs					
			alpha-BHC	0.02	--	ND	0.02	--
			beta-BHC	0.04	--	ND	0.2	--
			delta-BHC	0.02	--	ND	--	--
			gamma-BHC (Lindane)	0.03	--	ND	0.2	--
			Heptachlor	0.02	--	ND	0.4	--
			Aldrin	0.04	--	ND	0.04	--
			Heptachlor epoxide	0.05	--	ND	0.2	--
			Endosulfan I	0.04	--	ND	0.4	--
			Dieldrin	0.03	--	ND	0.03	--
			4,4'-DDE	0.04	--	ND	0.1	--
			Endrin	0.04	--	ND	2.0	--
			Endosulfan II	0.04	--	ND	0.4	--
			4,4'-DDD	0.04	--	ND	0.1	--
			Endosulfan sulfate	0.08	--	ND	0.4	--
			4,4'-DDT	0.04	--	ND	0.1	--
			Endrin aldehyde	0.1	--	ND	--	--
			Chlordane	0.1	--	ND	0.5	--
			Toxaphene	1.0	--	ND	3.0	--
			Aroclor-1016	1.0	--	ND	--	--
			Aroclor-1221	1.0	--	ND	--	--
			Aroclor-1232	1.0	--	ND	--	--
			Aroclor-1242	1.0	--	ND	--	--
			Aroclor-1248	1.0	--	ND	--	--
			Aroclor-1254	1.0	--	ND	--	--
			Aroclor-1260	1.0	--	ND	--	--

SOURCE: SMITH TECHNOLOGY CORPORATION (PROJECT NO. 09-5004-08)

TABLE 3

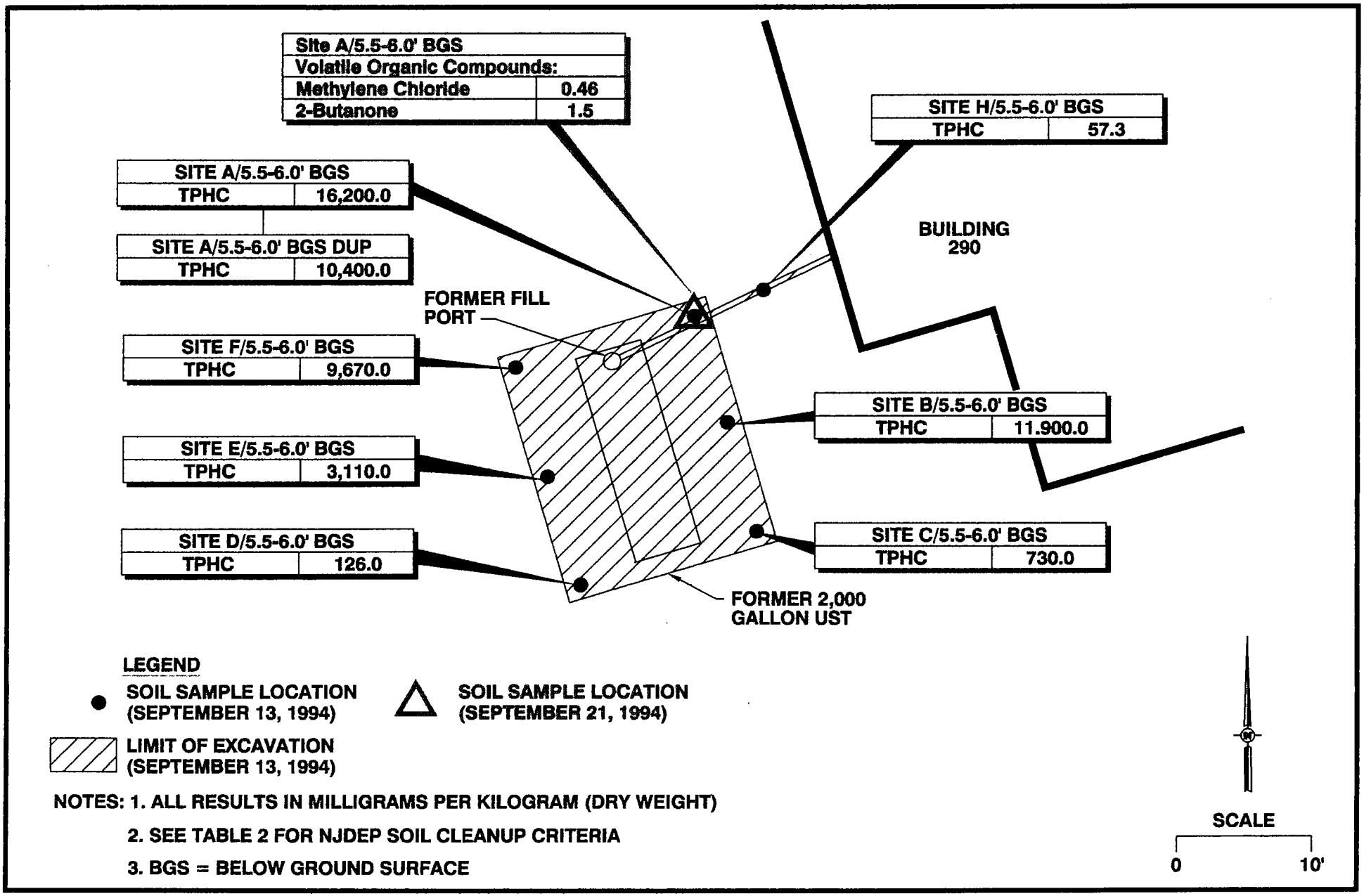
DATA ANALYSIS QUALIFIER DEFINITIONS
GROUNDWATER SAMPLING
FORT MONMOUTH, NEW JERSEY
PAGE 35 OF 35

PAGE 35 OF 35

--:	Not applicable / does not exceed criteria
J:	Indicates detected below sample quantitation limit
B:	Indicates also present in blank
ND:	Indicates compound not detected
NT	Not tested
NA:	Not available for this constituent
GWQS:	Groundwater Quality Standards

BCM ENGINEERS
A Division of

U.S. Army
Department of Public Works
Fort Monmouth, New Jersey



Project No. 09-5004-07

Figure 3
Building 290
Soil Sampling Results

APPENDIX A

NJDEP BUST CLOSURE APPROVAL

UNDERGROUND STORAGE TANK SYSTEM
CLOSURE APPROVAL

NEW JERSEY DEPARTMENT OF ENVIRONMENTAL PROTECTION AND ENERGY

DIVISION OF RESPONSIBLE PARTY SITE REMEDIATION
BUREAU OF UNDERGROUND STORAGE TANKS
CN-029, TRENTON, NJ 08625-0029

TMS #	UST #
C-93-3179	0081533

US Army
BLDG. 290
Ft. Monmouth, NJ

Monmouth

THE ABOVE LISTED FACILITY IS HEREBY GRANTED APPROVAL TO PERFORM THE FOLLOWING ACTIVITY IN ACCORDANCE WITH N.J.A.C. 7:14B-1 et. seq.:

Removal of: one 2,000 gallon #2 diesel UST(s) and appurtenant piping.
SITE ASSESSMENT: Soil samples will be taken every five (5) feet along the center line of each tank and one (1) soil sample for every 15 feet along all associated piping. Two (2) additional samples will be taken from around the tank and biased to the areas of highest field screened readings. Samples will be analyzed for TPHC. If sample results are greater than 1,000ppm than 25% of the samples will be analyzed for VO+10.

ON-SITE MANAGER: C. Appleby TELEPHONE: 908-532-1475

OWNER: TELEPHONE:

EFFECTIVE DATE: AUG 26 1993

THIS FORM MUST BE DISPLAYED AT THE SITE DURING THE APPROVED ACTIVITY AND MUST BE MADE AVAILABLE FOR INSPECTION AT ALL TIMES.

KEVIN F. KRATINA, BUREAU CHIEF
BUREAU OF UNDERGROUND STORAGE TANKS

APPENDIX B

CERTIFICATIONS

Let's preserve our earth

State of New Jersey
Department of Environmental Protection and Energy
Division of Responsible Party Site Remediation
CN 028
Trenton, NJ 08625-0029

ATTN: UST Program
(609) 984-3156

For State Use Only	
Date Rec'd.	
Auth.	
Routing	
UST NO.	

STANDARD REPORTING FORM
for reporting activities at an UST facility:

____ General Facility Information Changes	____ Sale or Transfer
_X_ Closure (Abandonment or Removal)	____ Substantial Modification
____ Temporary Closure	____ Financial Responsibility
____ Change in Service	____ Address Change Only

Check ONLY One Type of Activity – Complete Form For That Activity

(More than one tank can be listed per activity)

*** * * NOTE * * * ALL NEW tank installations at existing registered facilities must submit a Registration Questionnaire for the new tanks.**

Answer questions 1 through 5 and others as applicable.

1. Company name and address (as it appears on registration questionnaire): U.S. ARMY Fort Monmouth
DEH Bldg. 167, DEPT OF PUBLIC WORK
Fort Monmouth NJ 07703
ATTN: DINKER DESAI

2. Facility name and location (if different from above): _____

3. Contact person for this activity: DINKER DESAI
Telephone Number: (908) 532-1475

4. The identification number of the affected tank as it appears in Question Number 12 on the Registration Questionnaire: BLDG 290 C4

5. Registration Number (if known): UST - 0081533 - C

6. For GENERAL FACILITY INFORMATION changes (address, telephone, contact person, etc. – supply NEW information only):

   a. Facility name: _____
   b. Facility location: _____
   c. Owner's mailing address: _____
   _____
   _____ NJ _____
   _____
   d. Block: _____ Lot: _____
   e. Contact person (facility operator): _____
   f. Contact telephone number: (_____) _____ – _____
   g. Other (Specify): _____

(OVER)

7. For CLOSURE (abandonment or oval – check all that apply):
   a. ☐ Abandonment Date: ____/____/____ Case No: _____
      Attach the necessary implementation schedule (3 copies) and all documentation needed for abandonment per N.J.A.C. 7:14B-9.1 (d).
   b. ☒ Removal Date: 9/12/94 Case No. TMS # C-93-3179
      Attach the necessary implementation schedule (3 copies). 9-4-9-13-1303-57

8. For CHANGES IN HAZARDOUS SUBSTANCES STORED (check all that apply):
   a. ☐ Temporary Closure (12 month maximum time – see N.J.A.C. 7:14B-9.1(b)). Remove all hazardous substances; leave tank in place.
   b. ☐ Change in service from a regulated substance to a non-regulated substance. Tank must be cleaned and site assessment performed per N.J.A.C. 7:14B-9.1(e).
   c. ☐ Changes in service from one regulated hazardous substance to another regulated hazardous substance.
      Tank No. _____ Old _____ New _____
      Tank No. _____ Old _____ New _____
      Tank No. _____ Old _____ New _____
      (Attach additional sheets if more space is needed)

9. For TRANSFER OF OWNERSHIP: Effective Date: ____/____/____
   a. New Owner (operator) _____
   b. New Facility Name _____
      _____
      _____ NJ _____
      _____
      County
   c. Closing Attorney _____ Tele: (____) _____ - _____

10. For SUBSTANTIAL MODIFICATIONS (to include any retrofitted activity – e.g. the addition of spill/overfill protection, monitoring systems, cathodic protection, etc.):
   a. Type of Modification _____ Date: ____/____/____
   b. * NOTE * Substantial modifications require a permit under N.J.A.C. 7:14B-10.

11. For changes in FINANCIAL RESPONSIBILITY to (check appropriate changes and attach copies of new information):
   a. Policy Type: ☐
   b. Policy Number: ☐
   c. Other: ☐
   d. Company/Carrier: ☐
   e. Expiration Date: ☐

   _____
   _____
   _____
   (Specify)

NOTE: ALL appropriate and applicable permits, licenses and certificates required by the above activity(ies) from any local, state and/or federal agencies must be obtained separately from this notification.

CERTIFICATION

This registration form shall be signed by the highest ranking individual at the facility with overall responsibility for that facility (N.J.A.C. 7:14B-2.3 (a) 1).

"I certify under penalty of law that the information provided in this document is true, accurate and complete. I am aware that there are significant civil and criminal penalties for submitting false, inaccurate or incomplete information, including fines and/or imprisonment."

Signature: James Ott

Name (print or type): JAMES OTT

Title: Acting Director, Directorate of Public Works Date: 9/20/94

SRF-1/90
(INI/MOD-2/92)

APPENDIX C

WASTE MANIFEST

UNDERGROUND STORAGE TANK REMOVAL (UST)

(Submit one form for each tank)

Building No. 290 **NJDEPE UST Reg. No.** 0081533 - 64

IJO No. 91-0148 **Date Tank Removed** 9/12/94

ITEM NO.	ITEM OF WORK	UNIT	UNIT PRICE	QUANTITY	TOTAL PRICE
01100-1.1	Rmv ID#27 soil to stockpile	TN	$14.50	75.17	$1,089.97
01100-1.2	Supply, fill & relocate 55 Gal containers to storage	CT	$47.50		$
01100-1.4	Rmv & dispose of #2 fuel mixed with water Manifest #:NJA	GL	$ 0.69	5415 20	$ 3,750.15
01100-1.5	Rmv & dispose of #2 fuel mixed with solvent Manifest #:NJA	GL	$ 4.50		$
01100-1.6	Rmv & dispose of diesel fuel	GL	$ 0.69		$
01100-1.7	Rmv & dispose of diesel fuel mixed with water Manifest #:NJA	GL	$ 0.69		$
02050-1 & 02050-4	Tank removal	GL	$ 0.975	2000	$ 1,950.00
02050-5.1	Sawcut blacktop *	TN	$27.50	16.48	$ 453.20
02050-5.2	Sawcut concrete *	TN	$29.50	20.69	$ 610.36
02050-5.3	Sawcut *reinforced concrete*	TN	$32.50		$
02222-1.1	Backfill cert. clean fill *	TN	$16.25	48.75	$ 792.19
02222-1.2	3/4" clean stone *	TN	$17.50	26.42	$ 462.35
02511-1.1	Concrete slab 4" thick	SY	$19.80		$
02511-1.2	Concrete slab 6" thick	SY	$21.80		$
02511-1.3	Concrete slab 8" thick	SY	$24.50		$
02511-1.4	6" Concrete curb	LF	$16.00		$
02551-1.1	6" Base course of 3/4" dirty blend stone	SY	$ 6.40		$
02551-1.2	4" stabilized base	SY	8.00		$
02551-1.3	2" top course	SY	$ 5.50		$
02935-1.1	4" top soil & sod	SY	$ 7.80		$
02935-1.2	4" top soil & hydroseed	SY	$ 5.40		$

* Supply certified weight tickets to Contracting Officer at time of request for payment. 9,108.22

I certify under penalty of law that tank decommissioning activities were performed in compliance with NJAC 7:14B-9.2(b)3. I am aware that there are significant penalties for submitting false, inaccurate, or incomplete information, including fines and/or imprisonment.

NAME (Print or Type): John Lonergan SIGNATURE: John J Lonergan

NJDEPE UST Closure Cert. #: 0003248 DATE: 9/26/94

COMPANY NAME: CUTE Inc
(Performer of Tank Decommissioning)

NJDEPE UST Closure Corp. Cert. #: 0200128

List of Abbreviations:

CT = 55 Gallon Container GL = Gallon TN = Tons
EA = Each SY = Square Yard

FREEHOLD CARTAGE, INC.

P.O. BOX 5010
FREEHOLD, NJ 07728-5010
PHONE: (908) 462-1001
FAX: (908) 308 0924

175 BARTOW MUN. AIRPORT
BARTOW, FL 33830
PHONE: (813) 533-4599
FAX: (813) 533-1613

108 MONAHAN AVENUE
DUNMORE, PA 18512
PHONE: (717) 342-7232
FAX: (717) 342-7367

350 PIGEON POINT RD.
NEW CASTLE, DE 19720
PHONE: (302) 658-2005
FAX: (302) 658-6229

MANIFEST

FCI EPA ID NO.:
NJD054126164

G 59303

GENERATOR NAME/ADDRESS	PHONE (AREA CODE)	GENERATOR EPA ID NO.
US Army Communication & Electronics Command Main Post Fort Monmouth NJ	(908) 532-6224	
	TRACTOR: 63 / TRAILER: 252	APPOINTMENT TIME:

FCI REP. LOADING (PRINT)	PROCEDURE	BOX SPOTTED	BOX REMOVED	TIME AT GENERATOR (MILITARY TIME ONLY)
David Smith	VAC	X	X	ARRIVAL TIME: 07:30 / DEPARTURE TIME: 09:00

COMMENTS OR DELAYS AT GENERATOR	EQUIPMENT USED

BROKER:

PO. NO#:

STATE MANIFEST NO.:

(X) HM		PROPER U.S. D.O.T. SHIPPING NAME	U.S. D.O.T. HAZARDOUS CLASS	NO. CONT.	CONT. TYPE	WASTE NO.	PACKING GROUP	NA/UN NO.	FORM	NET QUANTITY	UNIT MEASURE
	1	NON-HAZARDOUS WASTE-WATER	N/A	1	FT	N/A	N/A	N/A	L	5415	G
	2	From Bldg 290	LIST II	0081533-64		—					
	3										

SPECIAL HANDLING INSTRUCTIONS INCLUDING CONTAINER EXEMPTION (I.E., IDENTIFICATION SHIPMENT OF A NON-HAZARDOUS NATURE WHICH DOES NOT HAVE TO BE MANIFESTED).

OW-1905 Rel-16 NJDEPE 15939-25213

GENERATOR'S CERTIFICATION: This is to certify that the above named materials are properly classified, described, packaged, marked and labeled and are in proper condition for transportation according to the applicable regulations of the Department of Transportation, U.S. EPA and the State. The wastes described above were consigned to the Transported named. The Treatment, Storage or Disposal Facility can and will accept the shipment of hazardous waste, and has a valid permit to do so. I certify that the foregoing is true and correct to the best of my knowledge.

Payment to the contractor for waste removal does not constitute payment to the carrier and if the contractor does not pay the carrier, the generator is obligated to pay the agreed rate offered to the contractor.

GENERATOR'S SIGNATURE	PLEASE PRINT NAME/TITLE	DATE LOADED
X [signature] I HAVE READ THE ABOVE AND UNDERSTAND AND AGREE TO ALL OF ITS CONTENT.	[illegible] - SELFM-M-EV	09/28/94 MO. DAY YR.

TSDF NAME/ADDRESS	PHONE (AREA CODE)	TSDF EPA ID NO.
EI DUPONT COMPANY CHAMBERS WORKS RT 130 DEEPWATER NJ 08023	609 540-2773	NJD002385730
	TRACTOR: 63 / TRAILER: 252	APPOINTMENT TIME:

FCI REP. UNLOADING (PRINT)	PROCEDURE	BOX SPOTTED	BOX REMOVED	TIME AT TSDF (MILITARY TIME ONLY)
David Smith	Gravity	X	X	ARRIVAL TIME: : / DEPARTURE TIME: :

COMMENTS OR DELAYS AT TSDF	EQUIPMENT USED

TSDF SIGNATURE	PLEASE PRINT NAME/TITLE	DATE UNLOADED
X		09/28/94 MO. DAY YR.

AR H-0257 PC 944	ME ME-HWT-47 ME-WOT-47	MO H-1490	NOVA SCOTIA, CANADA NSC 000 147	QUEBEC, CANADA QC-6ML-047
CT CT-HW-307	MD HWH-167 91-OP-1765	ND WH-429	OH 333-HW	RI RI-535
DE DE-HW-203 DE-SW-203	MA MA-294	NH TNH-0047	OK 3358	TX 40705
IL SWH-1540	MN 61572	NJ S-2265 15939	ONTARIO, CANADA A 840943	WI 11602
		NY JA-113	PA PA-AH-0067	

White - FCI Original
Yellow - FCI Billing
Blue - FCI Office/Customer
Green - Retained by TSDF
Gold - Retained by Generator

G 59303

State of New Jersey
Department of Environmental Protection and Energy
Hazardous Waste Regulation Program
Manifest Section
CN 421, Trenton, NJ 08625-0421

Please type or print in block letters. (Form designed for use on elite (12-pitch) typewriter.) Form Approved. OMB No. 2050-0039. Expires 9-30-94

UNIFORM HAZARDOUS WASTE MANIFEST

1. Generator's US EPA ID No. | Manifest Document No. 097225 | 2. Page 1 of 1 | Information in the shaded areas is not required by Federal law.

3. Generator's Name and Mailing Address: US Army Communications Electronics Command
Main Post, c/o James Shirghio, Bldg 2504
ATTN: SELFM-DL-EM-HS, Fort Monmouth, NJ 07703
4. Generator's Phone (908) 532-6226

A. State Manifest Document Number: NJA 1907275
B. State Generator's ID-(Gen. Site Address): Main Post Ft Monmouth

5. Transporter 1 Company Name: Freehold Cartage, Inc. | 6. US EPA ID Number: NJD053612614
C. State Trans. ID-NJDEPE: 522605 | Decal No.: 55464

7. Transporter 2 Company Name | 8. US EPA ID Number
D. Transporter's Phone (908) 462-1001
E. State Trans. ID-NJDEPE | Decal No.

9. Designated Facility Name and Site Address: Lionetti Oil Recovery Co., Inc.
Runyon & Cheesequake Rds.
Old Bridge, NJ 08857 | 10. US EPA ID Number: NJD064644664
F. Transporter's Phone ( )
G. State Facility's ID
H. Facility's Phone (908) 721-0900

	11. US DOT Description (Including Proper Shipping Name, Hazard Class or Division, ID Number and Packing Group)	HM	12. Containers No.	Type	13. Total Quantity	14. Unit Wt/Vol	Waste No.
a.	Petroleum Oil RXXM.N.O.S. Class 3 (Petroleum Oil) Combustible Liquid UN 1270 PG III	X	001	TT	00452	G	X 172
b.	Petroleum Oil N.O.S. Class 3 (Petroleum Oil) Combustible Liquid UN 1270 PG III	X	001	TT	00020	G	X 172
c.	Petroleum Oil N.O.S. Class 3 (Petroleum Oil) Combustible Liquid UN 1270 PG III	X	001	TT	00885	G	X 172
d.	Petroleum Oil N.O.S. Class 3 (Petroleum Oil) Combustible Liquid UN 1270 PG III	X	001	TT	01265	G	X 172

J. Additional Descriptions for Materials Listed Above

				K. Handling Codes for Wastes Listed Above
a.	Petroleum Oil 70 % T,L Water 30 %	Petroleum Oil 70 % T,L Water 30 %		T04=Filtration T04=Filtration
b.	Petroleum Oil 70 % T,L Water 30 %	Petroleum Oil 70 % T,L Water 30 %		T04=Filtration T04=Filtration

15. Special Handling Instructions and Additional Information
NOT REGULATED BY EPA. REGULATED AS HAZARDOUS WASTE IN NJ
24 HOUR EMERGENCY RESPONSE PHONE: 201-427-2881
NJ DECAL# 55464
a) 008153-67 b) 008153-64 c) 008153-63 d) 008153-68

16. GENERATOR'S CERTIFICATION: I hereby declare that the contents of this consignment are fully and accurately described above by proper shipping name and are classified, packed, marked, and labeled, and are in all respects in proper condition for transport by highway according to applicable international and national government regulations.

If I am a large quantity generator, I certify that I have a program in place to reduce the volume and toxicity of waste generated to the degree I have determined to be economically practicable and that I have selected the practicable method of treatment, storage, or disposal currently available to me which minimizes the present and future threat to human health and the environment; OR, if I am a small quantity generator, I have made a good faith effort to minimize my waste generation and select the best waste management method that is available to me and that I can afford.

Printed/Typed Name	Signature	Month	Day	Year
D. A. Detzel		09	01	94

TRANSPORTER

17. Transporter 1 Acknowledgement of Receipt of Materials

Printed/Typed Name	Signature	Month	Day	Year
BILL BURR		09	01	94

18. Transporter 2 Acknowledgement of Receipt of Materials

Printed/Typed Name	Signature	Month	Day	Year

FACILITY

19. Discrepancy Indication Space

20. Facility Owner or Operator: Certification of receipt of hazardous materials covered by this manifest except as noted in item 19.

Printed/Typed Name	Signature	Month	Day	Year
UDAL TAGAN	U.S. Tagan	09	02	94

EPA Form 8700-22 (Rev. 9/88) Previous editions are obsolete.

SIGNATURE AND INFORMATION *MUST* BE LEGIBLE ON ALL COPIES

STAVOLA ASPHALT CO.-TIN. FALLS

OFFICE COPY

Stavola Asphalt Co., Inc.

ASPHALT PLANTS

CONTROL NO. B374334

Old Bergen Mill Rd.
Millstone, NJ
(908) 446-7700

Hamilton Rd.
Tinton Falls, NJ
(908) 542-2328

Yellowbrook Rd.
Farmingdale, NJ
(908) 938-2801

Ashphalt from Bldg 290

THESE AREAS MUST BE SIGNED PRIOR TO UNLOADING OF MATERIAL

NOTE ATTENTION DRIVER: CUSTOMER MUST SIGN BELOW AND RETURN YELLOW COPY TO SCALE

X Don Ellis
DRIVER'S SIGNATURE

RECEIVED & ACCEPTED BY:
X
CUSTOMER'S SIGNATURE

EXECUTIVE OFFICE
HAMILTON ROAD
TINTON FALLS, NJ
908 / 542-2328

• ASPHALT
CRUSHED STONE • SAND
• GRAVEL

ADDRESS REPLY TO
P.O. BOX 482
RED BANK, N.J. 07701

THIS COMPANY WILL NOT BE RESPONSIBLE FOR DAMAGE CAUSED BY VEHICLES DELIVERING MATERIALS OFF PUBLIC ROADS.

EXPLANATION OF DELIVERY CODES
1 - F.O.B.
2 - DELIVERED
3 - NET DELIVERED

DATE	09/12/94	CUST. NO.	08888	JOB NO.	13:32	TICKET NO.	374334

CUSTOMER: CLEAN UP ENIVERMENT

DELIVER TO ZONE: Blacktop

GROSS 3.00 60380
TARE 2.00 27420
NET 1.00 32960

TRUCKER	TRUCK NO.	DRIVER NO.	METHOD OF PAYMENT	DELIVERY CODE	ZONE
08888	0		CHARGE	1	16.48T

QUANTITY	PRODUCT CODE/DESCRIPTION	UNIT OF MEASURE	UNIT PRICE	EXTENDED	FREIGHT	SALES TAX	TOTAL
1.00	48 RAP INCOMING	LD.					

WAITING TIME OVER 20 MINUTES MUST BE SIGNED BELOW BY CUSTOMER

H/T WAIT TIME 0.00

TIME: ON JOB _____ TIME: OFF JOB _____

~~LOADS~~ ~~ACCU. TONS~~
2 2.00

X _____
(SIG.)

GRAND TOTAL

STAVOLA ASPHALT CO.-TIN. FALLS

OFFICE COPY

Stavola Asphalt Co., Inc.

ASPHALT PLANTS

CONTROL NO. B374716

Old Bergen Mill Rd.
Millstone, NJ
(908) 446-7700

Hamilton Rd.
Tinton Falls, NJ
(908) 542-2328

Yellowbrook Rd.
Farmingdale, NJ
(908) 938-2801

Concrete
292+290

THESE AREAS MUST BE SIGNED PRIOR TO UNLOADING OF MATERIAL

NOTE ATTENTION DRIVER:
CUSTOMER MUST SIGN BELOW
AND RETURN YELLOW COPY TO SCALE

X Don Ellis
DRIVER'S SIGNATURE

RECEIVED & ACCEPTED BY:
X
CUSTOMER'S SIGNATURE

EXECUTIVE OFFICE
HAMILTON ROAD
TINTON FALLS, NJ
908 / 542-2328

• ASPHALT
CRUSHED STONE • SAND
• GRAVEL

ADDRESS REPLY TO
P.O. BOX 482
RED BANK, N.J. 07701

THIS COMPANY WILL NOT BE RESPONSIBLE FOR DAMAGE CAUSED BY VEHICLES DELIVERING MATERIALS OFF PUBLIC ROADS.

EXPLANATION OF DELIVERY CODES
1 - F.O.B.
2 - DELIVERED
3 - NET DELIVERED

DATE	09/15/94	CUST. NO.	08888	JOB NO.	11:21	TICKET NO.	374716

CUSTOMER: CLEAN UP ENIVERMENT

DELIVER TO ZONE: Concrete

GROSS 3.00 68,800
TARE 2.00 27420
NET 1.00 20,69

TRUCKER	TRUCK NO.	DRIVER NO.	METHOD OF PAYMENT	DELIVERY CODE	ZONE
08888	0		CHARGE	1	

QUANTITY	PRODUCT CODE/DESCRIPTION	UNIT OF MEASURE	UNIT PRICE	EXTENDED	FREIGHT	SALES TAX	TOTAL
1.00	48 RAP INCOMING	LD.					

WAITING TIME OVER 20 MINUTES MUST BE SIGNED BELOW BY CUSTOMER

H/WAIT TIME 0.00

TIME ON JOB	TIME OFF JOB	LOADS	ACCU. TONS
		2	2.00

X
(SIGNATURE)

GRAND TOTAL

CALCULATION SHEET

Building No. 290 NJDEPE Reg. No. 0081533 - 64

Tank Size 2000 gal Tank Void 15 tons

CLEAN FILL

ITEM NO.	DESCRIPTION	QUANTITY	TICKET #
	Fill	22.86	18820
		20.93	18821
		19.96	18822

TOTAL 63.75

STONE

ITEM NO.	DESCRIPTION	QUANTITY	TICKET #
	STONE	21.55	941555
		4.87	946850

TOTAL 26.42

ID#27 soil to stockpile (63.75+26.42) - 15 = 75.17 tons

Chargeable clean fill 63.75 - 15 = 48.75

Chargeable stone 26.42

S.C.M.I. - BOUND BROOK

OFFICE COPY

CONTROL NO. A-941555

Stavola Construction Materials, Inc.

PLANT: CHIMNEY ROCK ROAD, BOUND BROOK, N.J. • 908/356-5700

ATTENTION DRIVER:

PLEASE SIGN BELOW & RETURN GREEN "OFFICE COPY" TO SCALE BEFORE DEPARTING.

NOTE

X. Henry
DRIVER'S SIGNATURE

CRUSHED STONE • SAND • GRAVEL

EXECUTIVE OFFICE
HAMILTON ROAD
TINTON FALLS, N.J.
908/542-2328

ADDRESS REPLY TO
P.O. BOX 482
RED BANK, N.J. 07701

THIS COMPANY WILL NOT BE RESPONSIBLE FOR DAMAGE CAUSED BY VEHICLES DELIVERING MATERIALS OFF PUBLIC ROADS.

EXPLANATION OF DELIVERY CODES
1 - F.O.B.
2 - DELIVERED
3 - NET DELIVERED

DATE	CUST. NO.	JOB NO.	TICKET NO.
06/23/94.	08888	14:53	941555

CUSTOMER
CLEANING UP THE ENVIORNMENT
103 GODWIN AVE.
P.O. BOX 237
MIDLAND PARK NJ 07432

DELIVER TO
ZONE:FT MONMOUTH
CHARLES WOOD AREA

GROSS 35.55
TARE 14.00
NET 21.55

USER	TRUCK NO.	DRIVER NO.	METHOD OF PAYMENT	DELIVERY CODE	ZONE
03498	6		CHARGE	2	030

QUANTITY	PRODUCT CODE/DESCRIPTION	UNIT OF MEASURE	UNIT PRICE	EXTENDED	FREIGHT	SALES TAX	TOTAL
21.55	20 3/4 QPS-DGA	T			4.35		

LOADS 4 ACCU. TONS 94.33

WAIT TIME

GRAND TOTAL

M.I. BOUND BROOK

CUSTOMER'S COPY

CONTROL NO. A-946850

Stavola Construction Materials, Inc.

PLANT: CHIMNEY ROCK ROAD, BOUND BROOK, N.J. • 908/356-8700

X _____ DRIVER'S SIGNATURE

X _____ CUSTOMER'S SIGNATURE

CRUSHED STONE • SAND
GRAVEL

EXECUTIVE OFFICE
HAMILTON ROAD
TINTON FALLS, N.J.
908/542-2926

ADDRESS REPLY TO
P.O. BOX 482
RED BANK, N.J. 07701

THIS COMPANY WILL NOT BE RESPONSIBLE FOR DAMAGE CAUSED BY VEHICLES DELIVERING MATERIALS OFF PUBLIC ROADS.

EXPLANATION OF DELIVERY CODES
1 - F.O.B.
2 - DELIVERED
3 - NET DELIVERED

	CUST. NO.		JOB NO.		TICKET NO.	
07/08/94		08888		07158		946850

TOMER			DELIVER TO			GROSS		33.87
LEANING UP THE ENVIORNMENT.			ZONE #1 MONMOUTH BLDG 290			TARE		14.00
93 BODWIN AVE.						NET		19.87
P.O. BOX 237								
MIDLAND PARK NJ 07432								
ER	TRUCK NO.	DRIVER NO.	METHOD OF PAYMENT			DELIVERY CODE		ZONE
14010	39		CHARGE			2		030
TITY	PRODUCT CODE/DESCRIPTION		UNIT OF MEASURE	UNIT PRICE	EXTENDED	FREIGHT	SALES TAX	TOTAL
19.87	11 1.1/2" CLEAN STO		T			4.35		
MENTS						WAIT TIME		
			LOADS 2	ACCU. TONS 40.82		GRAND TOTAL		4.35 min

BLDG 108 7.5 tons
BLDG 443 7.5 tons
[redacted] 4.87 tons ~~Left~~
Bldg 290

Joseph Scerbo Sand & Gravel Co.

1455 W. Park Ave., Wayside
Asbury Park, N.J. 07712
908-493-3333

256

18820

Order Date Sept. 19, 94

Name Bill A Trucking

Deliver Date / /

Address Fill

Delivered ☐ C.O.D. ☐

F.O.B./P.U. ☐ Charge ☑

Item(s)	Quantity / Measure (tons, lbs., yds., ea.)	Unit Price	Total
	G 70121	22.86	
	T 24400		
	N 45721		

Driver

Received

Sub Total

Delivery

N.J. Tax

Total

* Company not responsible for damage done off public roads. Color not guaranteed

Have gravel will travel!
since 1925

PUBLIC WEIGHT MASTER No. 25614

Joseph Sansano Sand & Gravel Co.

1453 W. Park Ave., Wayside
Asbury Park, N.J. 07712
908-493-3333

2 FC

18821

Order Date Sept, 19, 94

Name Big A Trucking

Address Fill

Deliver Date ___/___/___

Delivered ☐ C.O.D. ☐

F.O.B./P.U. ☐ Charge ☑

Item(s)	Quantity / Measure (tons, lbs., yds., ea.)	Unit Price	Total
	G 70260		
	T 28400	20.93 tons	
	N 41860		
		Sub Total	
		Delivery	
		N.J. Tax	
		Total	

Driver ___

Received ___

No. 85514 (stamp: Public ... Weichert ...)

* Company not responsible for damage done on public roads. Color not guaranteed!

Have gravel will travel!
since 1925

Joseph Sorrento Sand & Gravel Co.

1453 W. Park Ave., Wayside
Asbury Park, N.J. 07712
908-493-3333

296

18822

Order Date Sept. 18, 99

Name Big A Trucking

Deliver Date ___/___/___

Address

Delivered ☐ C.O.D. ☐

Fill

F.O.B./P.U. ☐ Charge ☑

Item(s)	Quantity / Measure (tons, lbs., yds., ea.)	Unit Price	Total
	G 68335		
	T 28400	19.96 tons	
	W 39925		

KRISTEN WEICHERT PUBLIC No. 25914 NEW JERSEY

Driver

Received

Sub Total	
Delivery	
N.J. Tax	
Total	

* Company not responsible for damage done off public roads. Color not guaranteed!

Have gravel will travel!
since 1925!

APPENDIX D

UST DISPOSAL CERTIFICATE

TINTON FALLS, NJ

MAILING ADDRESS: 0 ASBURY AVE. NEPTUNE, NJ 07753

FACILITY ID NO: 1336F SP01

RECEIPT DOCUMENT NUMBER

01413600

HAULER

MARP508937
MARPAL COMPANY
PO BOX 188
LINCROFT NJ 07738

VEHICLE

TARE WEIGHT
18.1000 (36200)
GROSS WEIGHT
20.4900 (40980)

58523

DATE	OPER	ENTRY TIME	DEP NO	PLATE NO	EXIT TIME
10/24/94	JJJ	09:21	2065ZZ	XX77PH	09:38

Scale A (165) Normal

QUANTITY	CLASS	DESCRIPTION	UNIT	UNIT PRICE	AMOUNT
2.3900	13	Bulky Waste MONMOUTH COUNTY EATONTOWN BOROUGH	Tons	95.70	228.72

*** Prepayment Balance Remaining: 68740.04 ***

TRANSPORTER'S SIGNATURE

DOCUMENT TOTAL 228.72

CUSTOMER COPY

1,000 GAL FIBERGLASS TANK FROM BLDG 1104

3,000 GAL FIBERGLASS TANK FROM BLDG 290

MONITORING WELL CERTIFICATION-FORM B-LOCATION CERTIFICATION

Name of Permittee: U.S. ARMY
Name of Facility: FORT MONMOUTH
Location: MONMOUTH COUNTY, NJ
Case NJPDES Number: 93-11-30-1246-27

LAND SURVEYOR'S CERTIFICATION

Well Permit Number:
This number must be permanently affixed to the well casing. 29-309961-

Longitude (to nearest second): West 74° 02' 48.14"

Latitude (to nearest second): North 40° 18' 56.27"

Elevation of Top of Inner Casing (cap off) (one-hundredth of a foot): 13.90

Elevation of ground level (1/100th ft.) 16.42

Source of elevation datum (benchmark, nail, etc.) and year. (If an alternate datum has been approved by the Department, identify here, assume datum of 100', and give approximated actual elevation.) Source: FM-106

☒ 1927 ☐ 1983

Elev.: 17.80

Owners Well Number (As shown on application or plans): BLDG 290 MW-1

Elevations are to be determined by double run, three wire leveling methods using balanced sights, commencing from a well marked and described point. This beginning point shall either be derived from Federal or State benchmarks if not more than 1000 feet from the site or from an alternate datum approved by the Department. Tolerances should meet third order standards, which are 0.05 ft x $(\text{mile})^{1/2}$. For sections less than 0.1 mile, let miles = 0.1.

AUTHENTICATION

I certify under penalty of law that I have personally examined and am familiar with the information submitted in this document and all attachments and that, based on my inquiry of those individuals immediately responsible for obtaining the information, I believe the submitted information is true, accurate and complete. I am aware that there are significant penalties for submitting false information including the possibility of fine and imprisonment.

Wayne W Burgett
PROFESSIONAL LAND SURVEYOR'S SIGNATURE

WAYNE W BURGETT
PROFESSIONAL LAND SURVEYOR'S NAME
(Please print or type)

SEAL

31654
PROFESSIONAL LAND SURVEYOR'S LICENSE #

DWR-138 M
12/91

N. Jersey Department of Environmental Protection and Energy
Bureau of Water Allocation

MONITORING WELL RECORD

Well Permit No. _____ - _____
Atlas Sheet Coordinates _____ : _____ : _____ ☐

OWNER IDENTIFICATION - Owner _____
Address _____
City _____ State _____ Zip Code _____

WELL LOCATION - If not the same as owner please give address. Owner's Well No. [illegible]
County _____ Municipality _____ Lot No. _____ Block No. _____
Address _____

TYPE OF WELL (as per Well Permit Categories) _____ Date well completed 2/15/94
Regulatory Program Requiring Well _____ Case I.D. # _____

CONSULTING FIRM/FIELD SUPERVISOR (if applicable) _____ Tele. # _____

WELL CONSTRUCTION

Total depth drilled 12' ft.
Well finished to 12' ft.
Borehole diameter:
Top _____ in.
Bottom 6 in.
Well was finished: ☑ above grade ☐ flush mounted
If finished above grade, casing height (stick up) above land surface 3' ft.
Was steel protective casing installed? ☑ Yes ☐ No

	Depth to Top (ft.) [From land surface]	Depth to Bottom (ft.)	Diameter (inches)	Type and Material
Inner Casing	0	7'	4"	PVC
Outer Casing (Not Protective Casing)				
Screen (Note slot size)	7'	12½'	4"	.30 slot PVC
Tail Piece				
Gravel Pack	1'	12½'		#2
Annular Seal/Grout	0	1'		Concrete
Method of Grouting	Pour			

Static water level after drilling 5' ft.
Water level was measured using tape
Well was developed for 1 hours at 10 gpm
Method of development pump
Was permanent pumping equipment installed? ☐ Yes ☑ No
Pump capacity _____ gpm
Pump type: _____
Drilling Method auger
Drilling Fluid _____ Type of Rig CME-80
Name of Driller [illegible]
Health and Safety Plan submitted? ☑ Yes ☐ No
Level of Protection used on site (circle one) None D C B A
N.J. License No. 4431
Name of Drilling Company _____

GEOLOGIC LOG (Copies of other geologic logs and/or geophysical logs should be attached.)

0'-7' Clay - brown clay
FILL
7'-12' brown sand - FILL
12'-12½' brown clay - FILL

I certify that I have drilled the above-referenced well in accordance with all well permit requirements and all applicable State rules and regulations.

Driller's Signature [signature] Date 2-17-94

COPIES: White & Green - DEPE Canary - Driller Pink - Owner Goldenrod - Health Dept.

FIELD LOG OF BORING

SHEET ____ OF ____

LOCATION OF BORING: USArmy Ft Monmouth. Bldg 290 mw1

292 | 290 • mw1 | 296

PROJECT: US Army Ft Monmouth | BORING NO: mw1

TOTAL DEPTH: 12.5'

JOB NO: SELF DW EV | LOGGED BY: V Swanson

PROJ. MGR.: J Capriotti | EDITED BY:

DRILLING CONTRACTOR: Tyree

DRILL RIG TYPE: Mobil B50

DRILLERS NAME: Mike Beck

SAMPLING METHODS: split spoon

HAMMER WT.: 140 lb | DROP: 30"

STARTED, TIME: 1030 | DATE: 7/15/94

COMPLETED, TIME: 1330 | DATE: 7/15/94

BORING DEPTH (ft):	12.5
CASING DEPTH (ft):	2
WATER DEPTH (ft):	2
TIME:	1330
DATE:	7/15/94

BACKFILLED, TIME: 1330 DATE: 7/15/94 BY: Tyree

SURFACE ELEV: | DATUM:

CONDITIONS:

SAMPLE DEPTH	SAMPLER TYPE	BLOWS / 6 IN.	INCHES DRIVEN	INCHES RECOVERED	SAMPLE CONDITION	DRILLING RATE (min./ft.)	PID READING (ppm)	ODOR (Y / N ?)	GRAPHIC WELL CONST.	DEPTH IN FEET	GRAPHIC LOG
0-2	SS	1	6	12			0	N	casing & vented gravel pack	1	SC
		2	6								
		2	6							2	
		3	6				0	N			
										3	
							0	N		4	
										5	
							0	N	Screen	6	
										7	
							0	N		8	
										9	
							0	N		10	
							0	N		12.5	

SAND, grey-brown, medium and CLAY, till, trace

hit water at 2ft did 0-2' bori. twice because of no recovery the first time due to hitting a small rock.

VOTIS Sample.

TYREE ENVIRONMENTAL TECHNOLOGIES

GEOLOG: Drilling Activity Log Form

CLIENT: Fort Monmouth
PROJECT / JOB No.:
LOCATION: Bldg # 290
BOREHOLE / WELL No.:
LOGGED BY: Vaughn Swanson

DRILLING COMPANY: Tyree
DRILLER: m Beck
HELPER / CREW: J Rider
DRILLING TYPE: B80
DRILLER APPROVAL BY:

DATE	TIME: START	TIME: END	DRILLING ACTIVITY CODE	TOOL DIAM. (IN.)	SAMPLING METHOD	FOOTAGE DRILLED	EXPLANATION
7/15/94	1030	1330	WLC	4	SSC	25	

ACTIVITIES - GENERAL

- SBY = STANDBY
- DNT = DOWNTIME
- WLI = WELL INSTALLATION
- DCN = DECONTAMINATION
- WDV = WELL DEVELOPMENT
- GRT = GROUTING
- SPR = SITE SET UP/CLEAN UP
- TST = TESTING
- MOB = MOBILIZATION
- WKB = WORK BREAK
- WTR = FILL WATER TRUCK
- OTH = OTHER

ACTIVITIES - DRILLING

- HSA = HOLLOW STEM AUGERING
- MDR = MUD ROTARY
- ARR = AIR ROTARY
- RTF = ROTARY PUSH
- CBT = CABLE TOOL
- REM = REAMING
- DCD = DRILL THROUGH CASING DRIVER
- DUR = DUAL ROTARY CASING DRIVER
- COR = CORING
- RVC = REVERSE CIRCULATION
- DHH = DOWNHOLE HAMMER
- DAW = DRIVE CASING & WASH
- OTH = OTHER

SAMPLING (TESTING) METHODS

- SS/5 = 5-FT SPLIT SPOON SAMPLING
- SS/C = CONTINUOUS SPLIT SPOON SAMPLING
- CS/? = CALIFORNIA SPLIT SPOON SAMPLING (REPLACE ? WITH SAMPLING INTERVAL)
- CTS = CONTINUOUS TUBE SAMPLING
- STS = SHELBY TUBE SAMPLING
- COR = CORING
- WTS = WATER SAMPLING
- PKT = PACKER TESTING
- BGP = BOREHOLE GEOPHYSICS
- PMT = PUMPING TESTING
- SGT = SLUG TESTING
- OTH = OTHER

SUPPORT SERVICES OR DRILLING/WELL SUPPLIES (TYPE & SIZE)	DATE	HOURS	QUANTITY	UNIT

APPENDIX F

SOIL ANALYTICAL DATA PACKAGE

Report of Analysis
U.S. Army, Fort Monmouth Environmental Laboratory
NJDEPE Certification # 13461

Client: U.S. Army
DPW, SELFM-PW-EV
Bldg. 167
Ft. Monmouth, NJ 07703

Lab. ID #: 1641.1-.8
Sample Rec'd: 09/13/94
Analysis Start: 09/14/94
Analysis Comp: 09/14/94

Analysis: 418.1 (TPH)
Matrix: Soil
Analyst: S. Hubbard
Ext. Meth: Sonc.

NJDEPE UST Reg.#: **0081533-64**
Closure #: **C93-3179**
DICAR #: **9-4-9-13-1503-57**
Location #: **Bldg. 290**

Lab ID.	Description	%Solid	Result (mg/Kg)	MDL
1641.1	Site A, Sidewall N. OVA= 30.	86	16200.	130
1641.2	Site B, Sidewall NE. OVA= 20.	88	11900.	130
1641.3	Site C, Sidewall SE. OVA= 10.	83	730.	6.6
1641.4	Site D, Sidewall SO. OVA= 3.	82	126.	6.6
1641.5	Site E, Sidewall SW. OVA= 14.	84	3110.	46.
1641.6	Site F, Sidewall SE. OVA= 16.	85	9670.	46.
1641.7	Site G, Dup OVA= 30.	85	10400.	46.
1641.8	Site H, Pipe OVA= ND	83	57.3	9.9
M. Bl.	Method Blank	100	ND	3.3

Notes: ND = Not Detected, MDL = Method Detection Limit
* = Silica Gel Added, NA = Not Applicable
1641.3dup= 98% 1641.3s= 61% 1641.3sd= 63% RPD= 3.2%
Cal Chk = 102%

Brian K. McKee
Laboratory Director

Report of Analysis
U.S. Army, Fort Monmouth Environmental Laboratory
NJDEPE Certification # 13461

Client: U.S. Army
DPW, SELFM-PW-EV
Bldg. 167
Ft. Monmouth, NJ 07703

Lab. ID #: 1641.1-.8
Sample Rec'd: 09/13/94
Analysis Start: 09/14/94
Analysis Comp: 09/14/94

Analysis: Munsel

Lab ID#	Soil Color
1641.1	5Y 3/2 Dark Olive Gray
1641.2	5Y 3/2 Dark Olive Gray
1641.3	5Y 3/2 Dark Olive Gray
1641.4	5Y 3/2 Dark Olive Gray
1641.5	5Y 2.5/1 Black
1641.6	5Y 2.5/1 Black
1641.7	5Y 2.5/1 Black
1641.8	5Y 4/3 Olive

Brian K. McKee
Laboratory Director

U.S. ARMY FORT MONMOUTH

P.O. #: P4B7 TPHC

Chain of Custody

Project #: C93-3179	Sampler: George / Cube	Date / Time: Sept 13 / 9-50	Analysis Parameters: TPH, B/Shift, mOCH4, H-new		Start:	
Customer: Dinker Dalal DPW-ENVIRO	Site Name: BLDG 290 08 1533-64 C93-3179	Dicem # 9-4-9-13 1503-57			Finish:	
Phone: x 21457					Preservation Method	

Lab Sample ID Number	Date	Time	Customer Sample Location/ID Number	Sample Matrix	# of Bottles	TPH						H-new	Remarks
1641.1	9/13	10-14	Site A (NE) Sidwall	Soil	1	x	x	x				30	
.2	"	10-09	Site B NE	"	"	x	x	x				20	< 4°C
.3	"	10-18	Site C (SE)	"	"	x	x	x				10	
.4	"	10-21	Site D (SO)	"	"	x	x	x				3	4H-new
.5	"	10-25	Site E (SW)	"	"	x	x	x				14	
.6	"	10-26	Site F (SE)	"	"	x	x	x				16	
.7	"	10-12	Site G (Dup)	"	"	x	x	x				30	
.8	"	10-05	Site H (Pile)	"	"	x	x	x				ND	
													SN# 270136

Relinquished By (signature)	Date / Time	Received By (signature)	Shipped By: Hand
Relinquished By (signature)	Date / Time: 9/13 / 14-30	Received for Lab by (signature): Sarah Hubbard	Date / Time: 9/13/94 / 1430

Note: A drawing depicting sample location should be attached or drawn on the reverse side of this chain of custody. Site map attached

SAI-ENV COC form 01 Page 1 of 2 Pages Rev. A Date: 02 Apr 93.

Enviornmental Laboratory

Certification Number 13461

DIRECTORATE OF ENGINEERING AND HOUSING
FORT MONMOUTH, NEW JERSEY 07703-5108

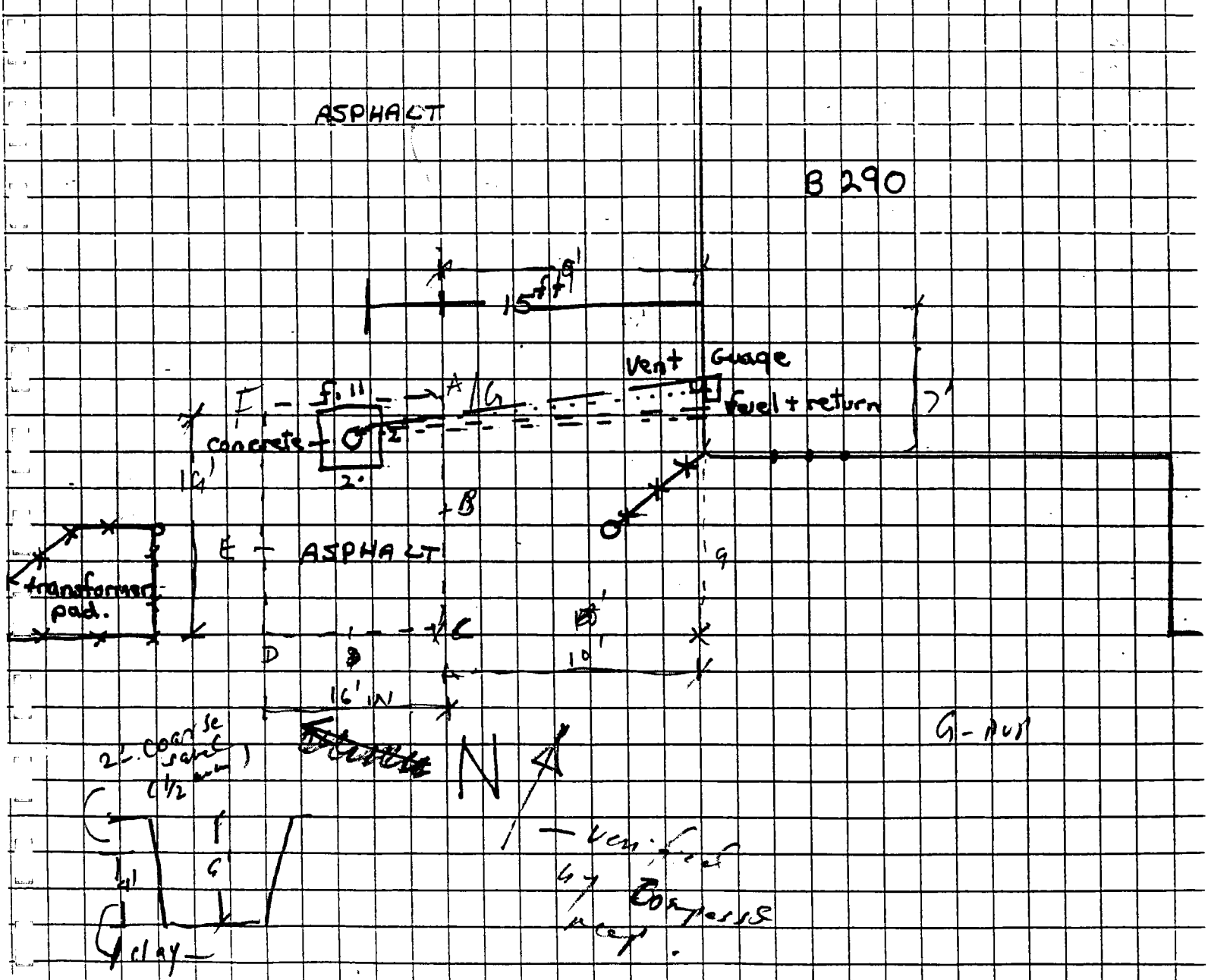
DATE: 11/23

Building Number: 290 Tank Number: 81533 10-64

Tank Size: 2000 gal Contents: #2 Fuel oil

Remarks:

1. All lines 15'
2.
3.



Sept 14, 1994 / Sarah O'Sullivan
1900

Blank 0 MV

40.75 59 MV

81.5 115 MV

163 234 MV

Method Blank 0 MV Building 292

1640.1 5 MV

1640.2 53 MV

1640.3 5 MV

1640.4 8 MV

1640.5 15 MV

1640.6 17 MV

1640.7 4 MV

1640.8 8 MV

1640.9 7 MV

40.75 Std 60 MV

Method Blank Building 290

1641.1 153 dil 13

1641.2 155 dil 13

1641.3 130 MV

1641.3 127 MV Dup

1641.3 191 MV Spk

1641.3 193 MV Dup Spk

1641.4 22 MV

1641.5 80 MV

195-6070-00

1641.54 22 MV

1641.55 80 MV

1641.6 168 dil 7

1641.7 180 dil 7

1641.8 10 MV

Method Blank Building 289

40.75 Standard 63 MV

1642.1 9 MV

1642.2 4 MV

1642.3 4 MV

1642.4 26 MV

1642.5 247 MV

1642.6 4 MV

1642.7 2 MV

1642.8 46 dil 7

40.75 Standard Ck 57 MV

Method Blank Bldg 618

1634.1 175 MV

1634.2 187 MV

1634.3 237 MV

1634.4 109 dil 7

1634.5 133 MV dil 7

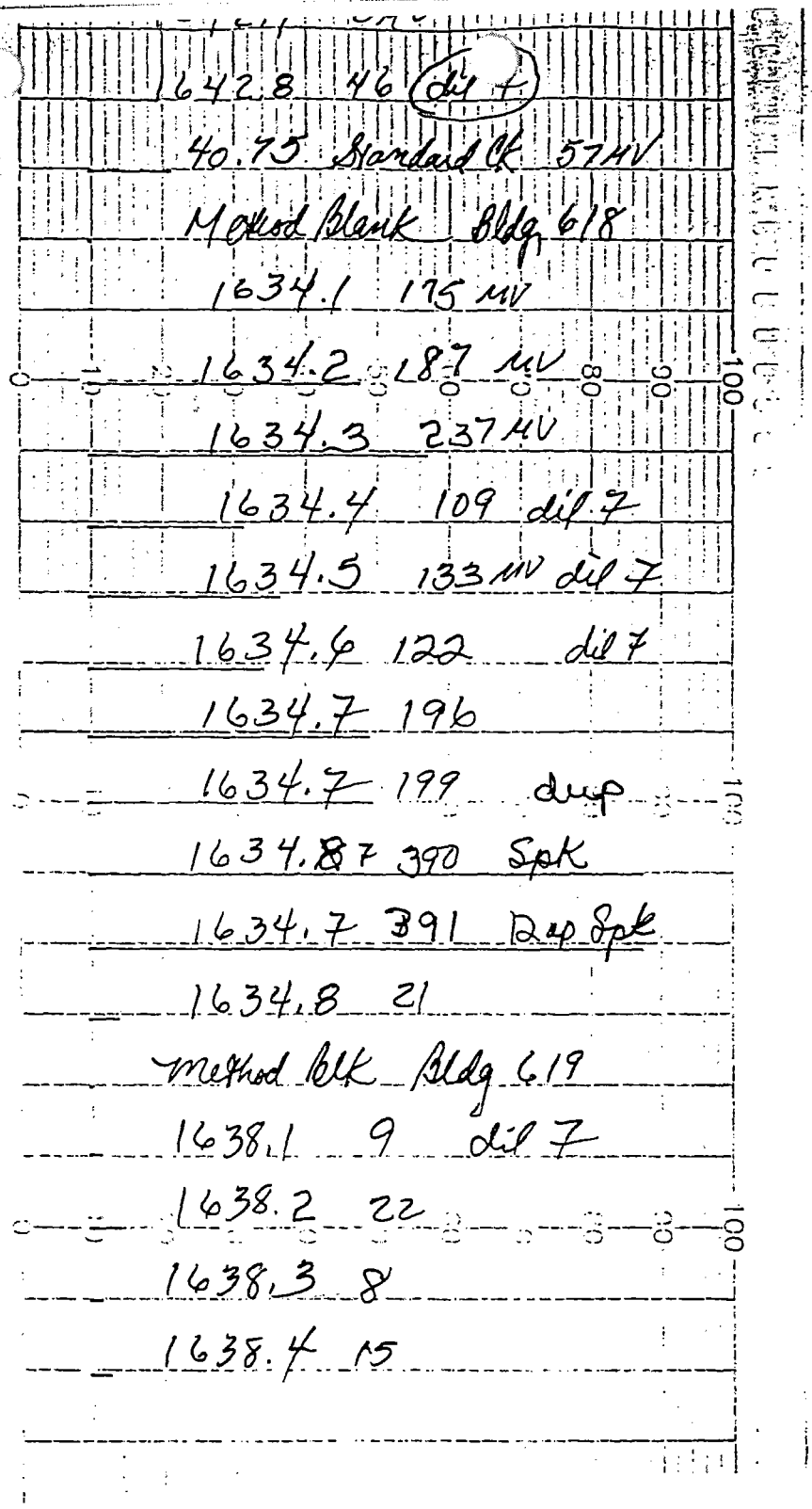
1634.6 122 dil 7

1634.7 196

1634.7 199 dup

1634.87 390 Spk

1634.7 391 Dup Spk



PHC Conformance/Non-conformance Summary Report

	No	Yes
1. Blank Contamination - If yes, list the sample and the corresponding concentrations in each blank	✓	
2. Matrix Spike/Matrix Sp Dup. Recoveries Meet Criteria (If not met, list the sample and corresponding recovery which falls outside the acceptable range)		✓
3. IR Spectra submitted for standards, blanks, & samples		✓
4. Chromatograms submitted for standards, blanks, and samples if GC fingerprinting was conducted.		N/A
5. Extraction holding time met. (If not met, list number of days exceeded for each sample)		✓
6. Analysis holding time met. (If not met,list number of days exceeded for each sample)		✓

Comments: _____

Laboratory Authentication Statement

I certify under penalty of law, where applicable, that this laboratory meets the Laboratory Performance Standards and Quality Control requirements specified in N.J.A.C. 7:18 and 40 CFR Part 136 for Water and Wastewater Analyses and SW 846 for Solid Waste Analysis. I have personally examined the information contained in this report, and to the best of my knowledge, I believe that the submitted information is true, accurate, complete, and meets the above referenced standards where applicable. I am aware that there are significant penalties for purposefully submitting falsified information, including the possibility of a fine and imprisonment.

Project #1641

Brian K. McKee
Laboratory Manager

princeton testing laboratory inc.

P.O. Box 3108
3490 U.S. Route 1
Princeton, NJ 08543-3108
(609) 452-9050
FAX (609) 452-0347

U.S. ARMY, FORT MONMOUTH
ATTN: SELFM-PW
Building 167
Fort Monmouth, New Jersey 07703-5108

Attn: Charles Appleby

Project # 94-8-11-1345-43
Building 482

JOB # 9404760-001

Laboratory Certification # 11118

Reviewed by: W. Alan Volk 11/4/94
W. Alan Volk

princeton testing laboratory inc.

P.O. Box 3108
3490 U.S. Route 1
Princeton, NJ 08543-3108
(609) 452-9050
FAX (609) 452-0347

TABLE OF CONTENTS

P404760

LABORATORY DELIVERABLES

THIS FORM MUST BE COMPLETED BY THE LABORATORY OR ENVIRONMENTAL CONSULTANT AND ACCOMPANY ALL DATA SUBMISSIONS

The following laboratory deliverables shall be included in the data submission. All deviations from the accepted methodology and procedures, or performance values outside acceptable ranges shall be summarized in the Non-Conformance Summary. The proposed "Technical Requirements for Site Remediation" rules, which appeared in the May 4, 1992 New Jersey Register, provides further details. The document shall be bound and paginated, contain a table of contents, and all pages shall be legible. Incomplete packages will be returned or held without review until the data package is completed.

It is recommended that the analytical results summary sheets listing all targeted and non-targeted compounds with the method detection limits be included in one section of the data package and in the main body of the report.

		Check if Complete
1.	Cover Page, Title Page listing Lab Certification #, facility name & address, & date of report	✓
2.	Table of Contents	✓
3.	Summary Sheets listing analytical results for all targeted and non-targeted compounds	✓
4.	Summary Table cross-referencing field ID #'s vs. Lab ID #'s	✓
5.	Document bound, paginated and legible	✓
6.	Chain of Custody	✓
7.	Methodology Summary	✓
8.	Laboratory Chronicle and Holding Time Check	✓
9.	Results submitted on a dry weight basis (if applicable)	✓
10.	Method Detection Limits	✓
11.	Lab certified by NJDEPE for parameters or appropriate category of parameters or a member of the USEPA CLP	✓
12.	Non-Conformance Summary	✓

Laboratory Manager or Environmental Consultant's Signature

4/4/94
Date

ORGANICS COMPLETE SDG FILE (CSF) INVENTORY SHEET

LABORATORY NAME Princeton Testing Laboratory
CITY/STATE Princeton, NJ.

CASE NO. 4760 SDG NO. 1544.1 SDG NOS. TO FOLLOW _____
_____ _____ SAS NO. _____

CONTRACT NO. Fort Monmouth
SOW NO. OLM01.8

All documents delivered in the complete SDG file must be original documents where possible. (REFERENCE EXHIBIT B, SECTION II and SECTION III.)

		PAGE NOs FROM	TO	CHECK LAB	EPA
1.	Inventory Sheet (Form DC-2) (Do not number)	___	___	✓	___
2.	SDG Case Narrative	___	___	✓	___
3.	SDG Cover Sheet/Traffic Report	___	___	✓	___
4.	Volatiles Data				
	a. QC Summary				
	System Monitoring Compound Summary (Form II VOA)	___	___	✓	___
	Matrix Spike/Matrix Spike Duplicate Summary (Form III VOA)	___	___	✓	___
	Method Blank Summary (Form IV VOA)	___	___	✓	___
	GC/MS Instrument Performance Check (Form V VOA)	___	___	✓	___
	Internal Standard Area and RT Summary (Form VIII VOA)	___	___	✓	___
	b. Sample Data	___	___		
	TCL Results - (Form I VOA)			✓	___
	Tentatively Identified Compounds (Form I VOA-TIC)			✓	___
	Reconstructed total ion chromatograms (RIC) for each sample			✓	___
	For each sample:				
	Raw spectra and background-subtracted mass spectra of target compounds identified			✓	___
	Quantitation reports			✓	___
	Mass spectra of all reported TICs with three best library matches			✓	___
	c. Standards Data (All Instruments)	___	___		
	Initial Calibration Data (Form VI VOA)			✓	___
	RICs and Quan Reports for all Standards			✓	___
	Continuing Calibration Data (Form VII VOA)			✓	___
	RICs and Quantitation Reports for all Standards			✓	___
	d. Raw QC Data				
	BFB	___	___	✓	___
	Blank Data	___	___	✓	___
	Matrix Spike/Matrix Spike Duplicate Data	___	___	✓	___

ORGANICS COMPLETE SDG FILE (CSF) INVENTORY SHEET (Cont.)

CASE NO. 4760 SDG NO. 1544.1 SDG NOS. TO FOLLOW _____
_____ _____ SAS NO. _____

	PAGE NOs FROM	TO	CHECK LAB	EPA
7. Miscellaneous Data				
Original preparation and analysis forms or copies of preparation and analysis logbook pages	___	___	___	___
Internal sample and sample extract transfer chain-of-custody records	___	___	___	___
Screening records	___	___	___	___
All instrument output, including strip charts from screening activities (describe or list)				
_____	___	___	___	___
_____	___	___	___	___
8. EPA Shipping/Receiving Documents				
Airbills (No. of shipments ___)	___	___	___	___
Chain-of-Custody Records	___	✓	___	___
Sample Tags	___	___	___	___
Sample Log-In Sheet (Lab & DCI)	___	___	___	___
Miscellaneous Shipping/Receiving Records (describe or list)				
_____	___	___	___	___
_____	___	___	___	___
9. Internal Lab Sample Transfer Records and Tracking Sheets (describe or list)				
_____	___	___	___	___
_____	___	___	___	___
10. Other Records (describe or list)				
Telephone Communication Log	___	___	___	___
_____	___	___	___	___
_____	___	___	___	___

11. Comments: _____

Completed by: (CLP Lab) _____ (Signature) W. A. Yolk QA/QC Coordinator (Printed Name/Title) 11/1/94 (Date)

Audited by: (EPA) _____ (Signature) _____ (Printed Name/Title) _____ (Date)

FORM DC-2-4

APPROVED SAMPLE ANALYSIS REQUEST

U.S. Army, Fort Monmouth N.J.
ATTN: SELFM-PW
Building 167
Fort Monmouth, New Jersey 07703-5108
Attention: Charles Appleby
Phone:(908) 532-6224 FAX:(908) 532-2367

Temp. Cust#: L9094
P.O. Number: E03-94U
Standard Tests

Project No.: 9404760-001M
Client Job#: 1644/1645
Date Received: 09/23/94
Analysis Due : 10/12/94

Number Of Samples : 3
Number Of Containers: 3

Approved By: Steven Burns

Reports: Custom Report Format

Sample I.D.'s	Code	Requested Analytical Services	Sampled
001 1645.1 Bldg 290 Site A Sidewall 9/21/94	VMS0B	Volatile Organics Library Search	09/21/94
	VMS0A	Volatile Organics, SW, SW-846 8240	
002 1644.1 Bldg 482 Site C-2 9/21/94 Sidewall SE	VMS0B	Volatile Organics Library Search	09/21/94
	VMS0A	Volatile Organics, SW, SW-846 8240	
003 1644.2 Fld Blk Bldgs 482/290 09/21/94	VMW0B	Volatile Organics Library Search	09/21/94
	VMW0A	Volatile Organics, WW, SW-846 8240	

Project Notes:

*** Results must be sent with ASCII disk. ***

Customer Notes:

Three Copies of Packages. See Data Mgmt for Details

Initials/Date

Received By Lab: _____
Reviewed By: _____
Q.A. Approved: _____

Printed By: Gene Dennison
Date: 11/01/94
Time: 12:31:29

940,166 001 M

U.S. ARMY FORT MONMOUTH

P.O. #: Princeton Lab 94-0551

Chain of Custody

Project #: 9-4-9-13-1503-57	Sampler: George / Cute	Date / Time 9/21/94 0900	Analysis Parameters: VOA +15	Start:
Customer: D. Desai Self - PW - EV	Site Name: Bldg. 290 UST 81533-64 TMS# C-93-3179 dlear 9-4-9-13 -1503-57			Finish:
Phone (908) 532-1475				Preservation Method

Lab Sample ID Number	Date/Time	Customer Sample Location/ID Number	Sample Matrix	# of Bottles	Remarks
1645.1	9/21/94 0945 ~~1005~~	Site A, Sidewall	soil	1	Shipped w/ ice
1644.2	9/21/94 0940	Field Blank	Aqua	1	

Relinquished By (signature)	Date / Time	Received By (signature)	Shipped By:
[signature]	9/21/94 1100	Sarah J. Hubbard	Hand
[signature]	9/23/94 1000	John Kennedy	9/23/94 1000

Note: A drawing depicting sample location should be attached or drawn on the reverse side of this chain of custody. Map Attached

Enviornmental Laboratory

N

Certification Number 13461

princeton testing laboratory inc.

P.O. Box 3108
3490 U.S. Route 1
Princeton, NJ 08543-3108
(609) 452-9050
FAX (609) 452-0347

LABORATORY CHRONICLE
ORGANIC ANALYSIS

Company: US Army, Fort Monmouth NJ Job #: 9404760-00/M

Date Received & Refrigerated: 09/23/94

EXTRACTION INFORMATION	ANALYSIS INFORMATION
Base-Neutral Extractables	Base-Neutral Extractables
___ ___/___/___	___ ___/___/___
___ ___/___/___	___ ___/___/___
___ ___/___/___	___ ___/___/___
Acid Extractables	Acid Extractables
___ ___/___/___	___ ___/___/___
___ ___/___/___	___ ___/___/___
___ ___/___/___	___ ___/___/___
Pesticides/ PCBs	Pesticides/ PCBs
___ ___/___/___	___ ___/___/___
___ ___/___/___	___ ___/___/___
PCBs only	PCBs only
___ ___/___/___	___ ___/___/___
___ ___/___/___	___ ___/___/___
Herbicides	Herbicides
___ ___/___/___	___ ___/___/___
___ ___/___/___	___ ___/___/___
Pesticides (EPTOX)	Pesticides (EPTOX)
___ ___/___/___	___ ___/___/___
Other: _____	Volatiles - 601/602
___ ___/___/___	___ ___/___/___
	___ ___/___/___
	Volatiles - 624/8240
	12 09/23/94
	___ 09/29/94
	___ 10/04/94
	Other: _____
	___ ___/___/___

Dept. Manager Review and Approval: [signature] 10/24/94

QC Supervisor Review and Approval: [signature] 11/4/94

princeton testing laboratory inc.

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3490 U.S. Route 1
Princeton, NJ 08543-3108
(609) 452-9050
FAX (609) 452-0347

METHODOLOGY SUMMARY

Laboratory: Princeton Testing Lab. Case Name: U.S. Army Ft. Monmouth

Location: Princeton, New Jersey Case Number: 9404760-001

VOLATILE ORGANIC ANALYSES:

EPA SW 846 8240

SEMIVOLATILE ORGANIC ANALYSES (ABN EXTRACTABLES):

PESTICIDES/PCBs and CHLORINATED HERBICIDES:

METALS ANALYSES:

TOTAL CYANIDE ANALYSES:

TOTAL PHENOL ANALYSES:

OTHER ANALYSES (SPECIFY):

NOTE: Only methods actually used in the performance of analyses for this data package may be entered on this form:
NJDEPE Form A-3 (9/91)

GCMS ANALYSIS NON CONFORMANCE SUMMARY

	NO	YES
1. GCMS TUNE SPECIFICATION.		
a. BFB Passed	____	✓
b. DFTPP Passed	____	____
2. GCMS TUNING FREQUENCY.		
a. Performed every 12 hours.	____	✓
b. Performed every 24 hours.	____	____
3. GCMS Calibration.		
a. Initial calibration performed w/i 30 days of sample analysis.	____	✓
b. Continuing calibration w/i 12 hours.	____	✓
c. Continuing calibration w/i 24 hours.	____	____
4. GCMS Calibration requirements.		
a. Calibration check compounds.	____	✓
b. System performance check compounds.	____	✓
5. Blank Contamination.		
a. VOA Fraction	10/04/94 Blank had Acetone below limits	
b. B/N Fraction		
c. Acid Fraction		
6. Surrogate Recoveries Within Limits.		
a. VOA Fraction	____	✓
b. B/N Fraction	____	____
c. Acid Fraction	____	____
7. Extraction Holding Time Met.	____	✓
8. Analysis Holding Time Met.		
a. VOA Fraction	____	✓
b. BNA Fraction	____	____

Comments:

Laboratory Manager [signature] Date 10/24/94

princeton testing laboratory inc.

P.O. Box 3108
3490 U.S. Route 1
Princeton, NJ 08543-3108
(609) 452-9050
FAX (609) 452-0347

October 21,1994.

U.S.Army, Fort Monmouth N.J
ATTN:SELFM-PW
Building 167
Fort Monmouth,New Jersey 07703-5108
Attention: Charles Appleby

Job Number: 9404760

CASE NARRATIVE

The following package contains analytical data pertaining to samples received by Princeton Testing Laboratory on 09/23/94. The samples were analyzed for volatile organics using SW-846, 8240 Methodologies.

VOLATILE ORGANICS

BLANKS: No contamination found in the blanks except 10/04 blank had acetone below Mdls.

SAMPLES: Methylene chloride and 2-butanone was found in samples 1645.1 and 1644.1.

SURROGATES: All surrogate recoveries were within QC limits.

MS/MSD: Sample 1639.1 Bldg 697 from the PTL Job# 9404683 was used for matrix spike and duplicate. All recoveries were within the QC limits.

If you have any further questions please do not hesitate to call me.

Khaja Eazazuddin.
GC/MS Supervisor.

S T A N D A R D T E S T L I S T

Test Name: Volatile Organics, SW, SW-846 8240
SPEC CODE: VMS0A

Lab Code: M
Manager : KE
Units : ug/kg

Short Name: VO.SW.8240
List Price:
Report Type:

Description

#	Compound	MDL	CAS#	KEY#	CLIENT KEY#
1	Chloromethane	10	00074873	C010	1454
2	Bromomethane	10	00074839	C015	1462
3	Vinyl chloride	10	00075014		1531
4	Chloroethane	10	00075003		1452
5	Methylene chloride	5	00075092		1172
6	Acetone	5	00067641		1498
7	Carbon disulfide	5	00075150		1646
8	1,1-Dichloroethene	5	00075354		1442
9	1,1-Dichloroethane	5	00075343		1580
10	1,2-Dichloroethene (Total)	5	00540590		1583
11	Chloroform	5	00067663		1078
12	1,2-Dichloroethane	5	00107062		1383
13	2-Butanone	5	00078933		1278
14	1,1,1-Trichloroethane	5	00071556		1068
15	Carbon tetrachloride	5	00056235		1384
16	Bromodichloromethane	5	00075274		1480
	1,1,2,2-Tetrachloroethane	5	00079345		1200
18	1,2-Dichloropropane	5	00078875		1368
19	trans-1,3-Dichloropropene	5	10061026		1364
20	Trichloroethene	5	00079016		1074
21	Dibromochloromethane	5	00124481		1390
22	1,1,2-Trichloroethane	5	00079005		1070
23	Benzene	5	00071432		1466
24	cis-1,3-Dichloropropene	5	10061015		1362
25	Bromoform	5	00075252		1124
26	2-Hexanone	5	00591786		1324
27	4-Methyl-2-Pentanone	5	00108101		1158
28	Tetrachloroethene	5	00127184		1096
29	Toluene	5	00108333		1426
30	Chlorobenzene	5	00108907		1564
31	Ethylbenzene	5	00100414		1394
32	Styrene	5	00100425		1120
33	Total Xylenes	5	01330207		1677
34	.				
35	.				
36	RECOVERY DATA	QC LIMITS			
37	.				
38	1,2-Dichloroethane-D4	70-121%			1577
39	Toluene-d8	84-138%	02037265		1579
40	4-Bromofluorobenzene	59-113%	00460004		1433

2B
VOLATILE SURROGATE SUMMARY

Lab Name:_Princeton Testing Lab. Contract:_US Army, Fort Monmouth.

Lab Code:_PTL_ Case No.:_4760_ SAS No.:_____ SDG No.:_____

Instrument ID: _INCOS-500___

SAMPLE NO.	(1,2-DCE)	(TOL-D8)	(4-BFB)
M.BLANK 9/23/94	111	104	101

(1,2-DCE)= 1,2-DICHLOROETHANE-d4 (76-114)
(TOL-d8)= TOLUENE-d8 (84-110)
(4-BFB) = 4-BROMOFLUOROBENZENE (86-115)

COMMENTS:

2B
VOLATILE SURROGATE SUMMARY

Lab Name:_Princeton Testing Lab. Contract:_US Army, Fort Monmouth.

Lab Code:_PTL_ Case No.:_4760_ SAS No.:_____ SDG No.:____

Instrument ID: _INCOS-500___

SAMPLE NO.	(1,2-DCE)	(TOL-D8)	(4-BFB)
1639.1 MS Bldg 697	71	93	77
1639.1 MSDBldg 697	72	88	79

(1,2-DCE)= 1,2-DICHLOROETHANE-d4 (70-121)
(TOL-d8)= TOLUENE-d8 (84-138)
(4-BFB) = 4-BROMOFLUOROBENZENE (59-113)

COMMENTS:

2B
VOLATILE SURROGATE SUMMARY

Lab Name:_Princeton Testing Lab. Contract:_US Army, Fort Monmouth.

Lab Code:_PTL_ Case No.:_4760_ SAS No.:_____ SDG No.:____

Instrument ID: _INCOS-500___

SAMPLE NO.	(1,2-DCE)	(TOL-D8)	(4-BFB)
M.BLANK 9/29/94	90	95	86
1645.1 9/21/94	80	90	86
1644.1 9/21/94	106	90	92

(1,2-DCE)= 1,2-DICHLOROETHANE-d4 (70-121)
(TOL-d8)= TOLUENE-d8 (84-138)
(4-BFB) = 4-BROMOFLUOROBENZENE (59-113)

COMMENTS:

page 1 of 1 FORM II VOA-2

2B
VOLATILE SURROGATE SUMMARY

Lab Name:_Princeton Testing Lab. Contract:_US Army, Fort Monmouth.

Lab Code:_PTL_ Case No.:_4760_ SAS No.:_____ SDG No.:_____

Instrument ID: _INCOS-500___

SAMPLE NO.	(1,2-DCE)	(TOL-D8)	(4-BFB)
M.BLANK 10/04/94	88	108	104
1644.2 09/21/94	96	105	92

(1,2-DCE)= 1,2-DICHLOROETHANE-d4 (76-114)
(TOL-d8)= TOLUENE-d8 (84-110)
(4-BFB) = 4-BROMOFLUOROBENZENE (86-115)

COMMENTS:

Princeton Testing Laboratory Inc.

P.O. Box 3108
3490 U.S. Route 1
Princeton, NJ 08543-3108
(609) 452-9050
(FAX) (609) 452-0347

Matrix Spike/Matrix Spike Duplicate Recovery Data
Test: Volatile Organics, Method 8240

Client: U.S. Army, Fort Monmouth N.J.

Project No.: 9404683-001

Lab Sample I.D.: 001

Client Sample I.D.: 1639.1 Bldg 697 MW 1 09/13/94

Analyst: Uma Chaudhary
Instrument: Incos 500 Volatiles
Units: ug/kg

QC Batch Number: 940923SV

COMPOUND	SPIKE ADDED	SAMPLE CONCENTRATION	MATRIX SPIKE CONCENTRATION	MS % REC	QC LIMITS REC
1,1-Dichloroethene	50	0	44.1	88.20	59-172
Trichloroethene	50	0	45.1	90.20	62-137
Benzene	50	0	44.3	88.60	66-142
Toluene	50	0	41.1	82.20	59-139
Chlorobenzene	50	0	46.6	93.20	60-133

COMPOUND	SPIKE ADDED	MSD CONCENTRATION	RPD	MSD % REC	QC LIMITS RPD
1,1-Dichloroethene	50	44.1	.00	88.20	0-22
Trichloroethene	50	48.7	7.68	97.40	0-24
Benzene	50	47.1	6.13	94.20	0-21
Toluene	50	43.3	5.21	86.60	0-21
Chlorobenzene	50	50.6	8.23	101.20	0-21

4A

VOLATILE METHOD BLANK SUMMARY

Lab Name: PTL,INC. Contract: US Army,Fort Monmouth.

Lab Code: PTL Case No.:4760-001SAS No.: xxxx SDG.No xxxx

Lab File ID: CBLK923... Lab Sample ID: LAB BLANK

Date Analyzed: 09/23/94 Time Analyzed: 12:08

Matrix (soil/water) SOIL... Level: (low/med) LOW

Instrument ID: FINN

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES, MS AND MSD:

	EPA SAMPLE NO.	LAB SAMPLE ID	LAB FILE ID	DATE ANALYZED
01	1639.1.MS...	4683-001-01.	.C8667...	09/23/94...
02	1639.1.MSD.	.4683-001-01.	.C8668...	09/23/94...
04	.....	.....	.....	.....
05	.....	.....	.....	.....

COMMENTS:

Princeton Testing Laboratory Inc.

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3490 U.S. Route 1
Princeton, NJ 08543-3108
(609) 452-9050
(FAX) (609) 452-1959

U.S. Army, Fort Monmouth N.J.
ATTN: SELFM-PW
Building 167
Fort Monmouth, New Jersey 07703-5108
Attention: Charles Appleby

Report Date: 10/20/94
Job Number: 9404760-001
Date Received: 09/23/94

Analysis: Volatile Organics, SW, SW-846 8240
Units: ug/kg

Parameters	*Sample I.D.: Blank 09/23/94*
Chloromethane	<10
Bromomethane	<10
Vinyl chloride	<10
Chloroethane	<10
Methylene chloride	<5.0
Acetone	<5.0
Carbon disulfide	<5.0
1,1-Dichloroethene	<5.0
1,1-Dichloroethane	<5.0
1,2-Dichloroethene (Total)	<5.0
Chloroform	<5.0
1,2-Dichloroethane	<5.0
2-Butanone	<5.0
1,1,1-Trichloroethane	<5.0
Carbon tetrachloride	<5.0
Bromodichloromethane	<5.0
1,1,2,2-Tetrachloroethane	<5.0
1,2-Dichloropropane	<5.0
trans-1,3-Dichloropropene	<5.0
Trichloroethene	<5.0
Dibromochloromethane	<5.0
1,1,2-Trichloroethane	<5.0
Benzene	<5.0
cis-1,3-Dichloropropene	<5.0
Bromoform	<5.0
2-Hexanone	<5.0
4-Methyl-2-Pentanone	<5.0
Tetrachloroethene	<5.0
Toluene	<5.0
Chlorobenzene	<5.0
Ethylbenzene	<5.0
Styrene	<5.0
Total Xylenes	<5.0

RECOVERY DATA	QC LIMITS	
1,2-Dichloroethane-d4 (Surrogate)	70-121%	111
Toluene-d8 (Surrogate)	84-138%	104
4-Bromofluorobenzene (Surrogate)	59-113%	101

EPA SAMPLE NO.
LAB BLK 9/23

1E
VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

Lab Name: Princeton Testing Lab US_ARMY,FORT_MONMOUTH.

Lab Code: PTL Case No.:4760-001SAS No.:XXX SDG No.:XXX

Matrix: (Soil/Water) Soil Lab Sample ID: LAB_BLK

Sample wt/vol: 5 (g/mL) g Lab File ID: CBLK923

Level: (low/med) LOW Date Received:

%Moisture: not dec. Date Analyzed: 09/23/94

GC Column: VOCOL ID: 0.53 mm Dilution Factor: 1

Soil Extract Vol: ul Soil Aliquot Vol: ul

Number TICs found: 0 CONCENTRATION UNITS:
(ug/L or ug/Kg) ug/Kg

#S	CAS NUMB	COMPOUND NAME	RT	EST. CONC.	SCAN

4A

VOLATILE METHOD BLANK SUMMARY

Lab Name: PTL,INC. Contract: US Army,Fort Monmouth.

Lab Code: PTL Case No.:4760-001SAS No.: xxxx SDG.No xxxx

Lab File ID: CBLK929A.. Lab Sample ID: LAB BLANK

Date Analyzed: 09/29/94 Time Analyzed: 10:26

Matrix (soil/water) SOIL.... Level: (low/med) LOW

Instrument ID: FINN

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES, MS AND MSD:

	EPA SAMPLE NO.	LAB SAMPLE ID	LAB FILE ID	DATE ANALYZED
01	1645.1.....	4760-001-01.	.C8694...	09/26/94...
02	1644.1.....	.4760-001-02.	.C8695...	09/26/94...
04	.....	.....	.....	.....
05	.....	.....	.....	.....

COMMENTS:

Princeton Testing Laboratory Inc.

P.O. Box 3108
3490 U.S. Route 1
Princeton, NJ 08543-3108
(609) 452-9050
(FAX) (609) 452-1959

U.S. Army, Fort Monmouth N.J.
ATTN: SELFM-PW
Building 167
Fort Monmouth, New Jersey 07703-5108
Attention: Charles Appleby

Report Date: 10/20/94
Job Number: 9404760-001
Date Received: 09/23/94

Analysis: Volatile Organics, SW, SW-846 8240
Units: ug/kg

Parameters	*Sample I.D.: Blank 09/29/94*
Chloromethane	<10
Bromomethane	<10
Vinyl chloride	<10
Chloroethane	<10
Methylene chloride	<5.0
Acetone	<5.0
Carbon disulfide	<5.0
1,1-Dichloroethene	<5.0
1,1-Dichloroethane	<5.0
1,2-Dichloroethene (Total)	<5.0
Chloroform	<5.0
1,2-Dichloroethane	<5.0
2-Butanone	<5.0
1,1,1-Trichloroethane	<5.0
Carbon tetrachloride	<5.0
Bromodichloromethane	<5.0
1,1,2,2-Tetrachloroethane	<5.0
1,2-Dichloropropane	<5.0
trans-1,3-Dichloropropene	<5.0
Trichloroethene	<5.0
Dibromochloromethane	<5.0
1,1,2-Trichloroethane	<5.0
Benzene	<5.0
cis-1,3-Dichloropropene	<5.0
Bromoform	<5.0
2-Hexanone	<5.0
4-Methyl-2-Pentanone	<5.0
Tetrachloroethene	<5.0
Toluene	<5.0
Chlorobenzene	<5.0
Ethylbenzene	<5.0
Styrene	<5.0
Total Xylenes	<5.0

RECOVERY DATA	QC LIMITS	
1,2-Dichloroethane-d4 (Surrogate)	70-121%	90
Toluene-d8 (Surrogate)	84-138%	95
4-Bromofluorobenzene (Surrogate)	59-113%	86

EPA SAMPLE NO.
LAB BLK 9/29

1E
VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

Lab Name: Princeton Testing Lab US_ARMY,FORT_MONMOUTH.

Lab Code: PTL Case No.:4760-001 SAS No.:XXX SDG No.:XXX

Matrix: (Soil/Water) Soil Lab Sample ID: LAB_BLK

Sample wt/vol: 5 (g/mL) g Lab File ID: CBLK929A

Level: (low/med) LOW Date Received:

%Moisture: not dec. Date Analyzed: 09/29/94

GC Column: VOCOL ID: 0.53 mm Dilution Factor: 1

Soil Extract Vol: ul Soil Aliquot Vol: ul

Number TICs found: 0 CONCENTRATION UNITS:
(ug/L or ug/Kg) ug/Kg

#S	CAS NUMB	COMPOUND NAME	RT	EST. CONC.	SCAN
1	1073-06-9	BENZENE,1-BROMO-3-FLUORO-	20:06	56	798

4A

VOLATILE METHOD BLANK SUMMARY

Lab Name: PTL,INC. Contract: US Army,Fort Monmouth.

Lab Code: PTL Case No.:4760-001SAS No.: xxxx SDG.No xxxx

Lab File ID: CBLK1004.. Lab Sample ID: LAB BLANK

Date Analyzed: 10/04/94 Time Analyzed: 13:33

Matrix (soil/water) WATER... Level: (low/med) LOW

Instrument ID: FINN

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES, MS AND MSD:

	EPA SAMPLE NO.	LAB SAMPLE ID	LAB FILE ID	DATE ANALYZED
01	1644.2.9/21	.4760-001-03.	.C8711...	10/04/94...
02				
03				

COMMENTS:

Princeton Testing Laboratory Inc.

P.O. Box 3108
3490 U.S. Route 1
Princeton, NJ 08543-3108
(609) 452-9050
(FAX) (609) 452-1959

U.S. Army, Fort Monmouth N.J.
ATTN: SELFM-PW
Building 167
Fort Monmouth, New Jersey 07703-5108
Attention: Charles Appleby

Report Date: 11/02/94
Job Number: 9404760-001
Date Received: 09/23/94
Client Job No.: 1644/1645
Page: 1

Analysis: Volatile Organics, WW, SW-846 8240
Units: ug/liter

Parameters	*Sample I.D.: Blank 10/04/94*
Chloromethane	<10
Bromomethane	<10
Vinyl chloride	<10
Chloroethane	<10
Methylene chloride	<5.0
Acetone	2.6 J
Carbon disulfide	<5.0
1,1-Dichloroethene	<5.0
1,1-Dichloroethane	<5.0
1,2-Dichloroethene (Total)	<5.0
Chloroform	<5.0
1,2-Dichloroethane	<5.0
2-Butanone	<5.0
1,1,1-Trichloroethane	<5.0
Carbon tetrachloride	<5.0
Bromodichloromethane	<5.0
1,1,2,2-Tetrachloroethane	<5.0
1,2-Dichloropropane	<5.0
trans-1,3-Dichloropropene	<5.0
Trichloroethene	<5.0
Dibromochloromethane	<5.0
1,1,2-Trichloroethane	<5.0
Benzene	<5.0
cis-1,3-Dichloropropene	<5.0
Bromoform	<5.0
2-Hexanone	<5.0
4-Methyl-2-Pentanone	<5.0
Tetrachloroethene	<5.0
Toluene	<5.0
Chlorobenzene	<5.0
Ethylbenzene	<5.0
Styrene	<5.0
Total Xylenes	<5.0

RECOVERY DATA	QC LIMITS	
1,2-Dichloroethane-d4 (Surrogate)	76-114%	88
Toluene-d8 (Surrogate)	88-110%	108
4-Bromofluorobenzene (Surrogate)	86-115%	104

J = Estimated Value Detected Below MDL

EPA SAMPLE NO.
LAB BLK 10/04

1E
VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

Lab Name: Princeton Testing Lab US_ARMY,FORT_MONMOUTH.

Lab Code: PTL **Case No.:** 4760-001 **SAS No.:** XXX **SDG No.:** XXX

Matrix: (Soil/Water) WATER

Lab Sample ID: LAB_BLK

Sample wt/vol: 5 (g/mL) mL

Lab File ID: CBLK1004

Level: (low/med) LOW

Date Received:

%Moisture: not dec.

Date Analyzed: 10/04/94

GC Column: VOCOL ID: 0.53 mm

Dilution Factor: 1

Soil Extract Vol: ul

Soil Aliquot Vol: ul

Number TICs found: 0

CONCENTRATION UNITS:
(ug/L or ug/Kg) ug/L

#S	CAS NUMB	COMPOUND NAME	RT	EST. CONC.	SCAN

5A

VOLATILE ORGANIC GC/MS TUNING AND MASS CALIBRATION - BROMOFLUOROBENZENE (BFB)

Lab Name: Princeton Testing Lab. Contract: US ARMY, FORT MONMOUTH

Lab Code: PTL Case No.: 4760 SAS No.: _____ SDG No.: _____

Lab File ID: BFB914 BFB Injection Date: 9/14/94

Instrument ID: FINN500V BFB Injection Time: 1200

Matrix:(soil/water) WATER Level:(low/med) Low Column:(pack/cap) Cap

m/e	ION ABUNDANCE CRITERIA	% RELATIVE ABUNDANCE
50	15.0 - 40.0% of mass 95	18.3
75	30.0 - 60.0% of mass 95	48.5
95	Base peak, 100% relative abundance	100.0
96	5.0 - 9.0% of mass 95	8.3
173	Less than 2.0% of mass 174	0.0
174	Greater than 50.0% of mass 95	86.4
175	5.0 - 9.0% of mass 174	7.4
176	Greater than 95.0%, but less than 101.0% of mass 174	99.3
177	5.0 - 9.0% of mass 176	6.9

1-Value is % mass 174 2-Value is % mass 176

THIS TUNE APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS, AND STANDARDS:

EPA SAMPLE NO.	LAB SAMPLE ID	LAB FILE ID	DATE ANALYZED	TIME ANALYZED
50 PPB STD	50 PPB STD	CV91450A	9/14/94	1943
10 PPB STD	10 PPB STD	CV91410A	9/14/94	2123
20 PPB STD	20 PPB STD	CV91420A	9/14/94	2034
100 PPB STD	100 PPB STD	CV914100B	9/14/94	1853
200 PPB STD	200 PPB STD	CV914200	9/14/94	1803

5A

VOLATILE ORGANIC GC/MS TUNING AND MASS CALIBRATION - BROMOFLUOROBENZENE (BFB)

Lab Name: Princeton Testing Lab. Contract: US ARMY, FORT MONMOUTH

Lab Code: PTL Case No.: 4760 SAS No.: _____ SDG No.: _____

Lab File ID: BFB923 BFB Injection Date: 9/23/94

Instrument ID: FINN500V BFB Injection Time: 1055

Matrix:(soil/water)SOIL Level:(low/med)Low Column:(pack/cap)Cap

m/e	ION ABUNDANCE CRITERIA	% RELATIVE ABUNDANCE
50	15.0 - 40.0% of mass 95	20.7
75	30.0 - 60.0% of mass 95	53.9
95	Base peak, 100% relative abundance	100.0
96	5.0 - 9.0% of mass 95	7.3
173	Less than 2.0% of mass 174	0.0
174	Greater than 50.0% of mass 95	52.7
175	5.0 - 9.0% of mass 174	7.2
176	Greater than 95.0%, but less than 101.0% of mass 174	99.3
177	5.0 - 9.0% of mass 176	7.0

1-Value is % mass 174 2-Value is % mass 176

THIS TUNE APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS, AND STANDARDS:

EPA SAMPLE NO.	LAB SAMPLE ID	LAB FILE ID	DATE ANALYZED	TIME ANALYZED
50_PPB STD	50 PPB STD	CV0923	9/23/94	1107
LAB_BLANK	M.BLK	CBLK923	9/23/94	1208
1639.1 MS	4683-001-01	C8667	9/23/94	1633
1639.1_MSD	4683-001-01	C8668	9/23/94	1723

5A

VOLATILE ORGANIC GC/MS TUNING AND MASS CALIBRATION - BROMOFLUOROBENZENE (BFB)

Lab Name: Princeton Testing Lab. Contract: US ARMY, FORT MONMOUTH

Lab Code: PTL Case No.: 4760 SAS No.: _____ SDG No.: _____

Lab File ID: BFB929 BFB Injection Date: 9/29/94

Instrument ID: FINN500V BFB Injection Time: 08:45

Matrix:(soil/water) SOIL Level:(low/med) Low Column:(pack/cap) Cap

m/e	ION ABUNDANCE CRITERIA	% RELATIVE ABUNDANCE
50	15.0 - 40.0% of mass 95	19.9
75	30.0 - 60.0% of mass 95	52.4
95	Base peak, 100% relative abundance	100.0
96	5.0 - 9.0% of mass 95	7.1
173	Less than 2.0% of mass 174	0.0
174	Greater than 50.0% of mass 95	65.2
175	5.0 - 9.0% of mass 174	7.1
176	Greater than 95.0%, but less than 101.0% of mass 174	97.8
177	5.0 - 9.0% of mass 176	6.8

1-Value is % mass 174 2-Value is % mass 176

THIS TUNE APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS, AND STANDARDS:

EPA SAMPLE NO.	LAB SAMPLE ID	LAB FILE ID	DATE ANALYZED	TIME ANALYZED
50_PPB STD	50 PPB STD	CV0929	9/29/94	0856
LAB_BLANK	M.BLK	CBLK929A	9/29/94	1026
1645.1	4760-001-01	C8694	9/29/94	1746
1644.1	4760-001-02	C8696	9/29/94	1928

5A

VOLATILE ORGANIC GC/MS TUNING AND MASS CALIBRATION - BROMOFLUOROBENZENE (BFB)

Lab Name: Princeton Testing Lab. Contract: US ARMY, FORT MONMOUTH

Lab Code: PTL Case No.: 4760 SAS No.: XXX SDG No.: XXX

Lab File ID: BFB1004B BFB Injection Date: 10/04/94

Instrument ID: FINN500V BFB Injection Time: 11:42

Matrix: (soil/water) WATER Level: (low/med) Low Column: (pack/cap) Cap

m/e	ION ABUNDANCE CRITERIA	% RELATIVE ABUNDANCE
50	15.0 - 40.0% of mass 95	21.3
75	30.0 - 60.0% of mass 95	52.7
95	Base peak, 100% relative abundance	100.0
96	5.0 - 9.0% of mass 95	7.3
173	Less than 2.0% of mass 174	0.0
174	Greater than 50.0% of mass 95	57.7
175	5.0 - 9.0% of mass 174	7.6
176	Greater than 95.0%, but less than 101.0% of mass 174	95.3
177	5.0 - 9.0% of mass 176	7.4

1-Value is % mass 174 2-Value is % mass 176

THIS TUNE APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS, AND STANDARDS:

EPA SAMPLE NO.	LAB SAMPLE ID	LAB FILE ID	DATE ANALYZED	TIME ANALYZED
50 PPB STD	50 PPB STD	CV0104	10/04/94	1234
LAB BLANK	M.BLK	CBLK1004	10/04/94	1333
1644.2	4760-001-03	C8711	10/04/94	1421

8A

VOLATILE INTERNAL STANDARD AREA AND RT SUMMARY

Lab Name: Princeton Testing Lab. Contract: US_Army,Fort_monmouth.___.
Case No.: 4760_ Lab File ID (standard):CV0923_____ Instrument ID:FINN
Date Analyzed: 09/23/94____ Time Analyzed: 11:07_____

		IS1(BCM)	RT	IS2(DFB)	RT	IS3(CB)	RT
	12hr. STD __	12697_____	07:35	158172_____	09:52	106406_____	22:18_
	Upper Limit_	25394_____	07:85	316344_____	10:02	21281_____	22:68_
	Lower Limit	6348_____	06:85	79086_____	09:02	53203_____	21:68_
	EPA Sample.						
1	M.BLANK__9/23	12703_____	07:29	121792_____	09:48	88184_____	22:01_
2	1639.1_MS____	8847_____	07:33	101972_____	09:49	74708_____	22:02_
3	1639.1_MSD___	8993_____	07:33	96476_____	09:49	72942_____	22:02_

IS1=BROMOCHLOROMETHANE
IS2=1,4-DIFLUOROBENZENE
IS3=CHLOROBENZENE-D5

Area Upper Limit = +100% of Internal Standard area.
Area Lower Limit = - 50% of Internal Standard area.
RT Upper Limit = +0.50 minutes of Internal Standard RT.
RT Lower Limit = -0.50 minutes of Internal Standard RT.

* Values outside of QC limits. FORM VIII V-1 PAGE 1 of 1.

8A

VOLATILE INTERNAL STANDARD AREA AND RT SUMMARY

Lab Name: Princeton Testing Lab. Contract: US_Army,Fort_monmouth.___.
Case No.: 4760_ Lab File ID (standard):CV0929_____ Instrument ID:FINN
Date Analyzed: 09/29/94____ Time Analyzed: 08:56_____

	IS1(BCM)	RT	IS2(DFB)	RT	IS3(CB)	RT
12hr. STD __	17120_____	07:30	142012_____	09:46	99603_____	21:58_
Upper Limit_	34240_____	07:80	284024_____	09:96	199206_____	22:08_
Lower Limit	8560_____	06:80	71006_____	08:96	49801_____	21:08_
EPA Sample.						
1 M.BLANK__9/29	10486_____	07:33	121433_____	09:48	89390_____	21:59_
2 1645.1_____	10152_____	07:30	68901_____	09:46	56616_____	22:01_
3 1644.1_____	8235_____	07:32	144726_____	09:48	114280_____	21:59_

IS1=BROMOCHLOROMETHANE
IS2=1,4-DIFLUOROBENZENE
IS3=CHLOROBENZENE-D5

Area Upper Limit = +100% of Internal Standard area.
Area Lower Limit = - 50% of Internal Standard area.
RT Upper Limit = +0.50 minutes of Internal Standard RT.
RT Lower Limit = -0.50 minutes of Internal Standard RT.

* Values outside of QC limits. FORM VIII V-1 PAGE 1 of 1.

8A

VOLATILE INTERNAL STANDARD AREA AND RT SUMMARY

Lab Name: Princeton Testing Lab. Contract: US_Army,Fort_monmouth.___.
Case No.: 4760_ Lab File ID (standard):CV0104_____ Instrument ID:FINN
Date Analyzed: 10/04/94____ Time Analyzed: 12:34_____

	IS1(BCM)	RT	IS2(DFB)	RT	IS3(CB)	RT
12hr. STD __	15702_____	07:22	113786_____	09:38	82666_____	21:48_
Upper Limit_	31404_____	07:72	227572_____	09:88	165332_____	21:98_
Lower Limit	7851_____	06:72	56893_____	08:88	41333_____	20:98_
EPA Sample.						
1 M.BLANK_10/04	10934_____	07:24	102320_____	09:38	73456_____	21:53_
2 1644.2_____	11875_____	07:22	87995_____	09:40	47327_____	21:53_

IS1=BROMOCHLOROMETHANE
IS2=1,4-DIFLUOROBENZENE
IS3=CHLOROBENZENE-D5

Area Upper Limit = +100% of Internal Standard area.
Area Lower Limit = - 50% of Internal Standard area.
RT Upper Limit = +0.50 minutes of Internal Standard RT.
RT Lower Limit = -0.50 minutes of Internal Standard RT.

* Values outside of QC limits. FORM VIII V-1 PAGE 1 of 1.

Princeton Testing Laboratory Inc.

P.O. Box 3108
3490 U.S. Route 1
Princeton, NJ 08543-3108
(609) 452-9050
(FAX) (609) 452-1959

U.S. Army, Fort Monmouth N.J.
ATTN: SELFM-PW
Building 167
Fort Monmouth, New Jersey 07703-5108
Attention: Charles Appleby

Report Date: 11/01/94
Job Number: 9404760-001
Date Received: 09/23/94
Client Job No.: 1644/1645

Analysis: Volatile Organics, SW, SW-846 8240
Units: ug/kg

Parameters	*Sample I.D.: 1645.1 Bldg 290 Site A Sidewall 9/21/94*
Chloromethane	<1200
Bromomethane	<1200
Vinyl chloride	<1200
Chloroethane	<1200
Methylene chloride	460 J
Acetone	<620
Carbon disulfide	<620
1,1-Dichloroethene	<620
1,1-Dichloroethane	<620
1,2-Dichloroethene (Total)	<620
Chloroform	<620
1,2-Dichloroethane	<620
2-Butanone	1500
1,1,1-Trichloroethane	<620
Carbon tetrachloride	<620
Bromodichloromethane	<620
1,1,2,2-Tetrachloroethane	<620
1,2-Dichloropropane	<620
trans-1,3-Dichloropropene	<620
Trichloroethene	<620
Dibromochloromethane	<620
1,1,2-Trichloroethane	<620
Benzene	<620
cis-1,3-Dichloropropene	<620
Bromoform	<620
2-Hexanone	<620
4-Methyl-2-Pentanone	<620
Tetrachloroethene	<620
Toluene	<620
Chlorobenzene	<620
Ethylbenzene	<620
Styrene	<620
Total Xylenes	<620

RECOVERY DATA	QC LIMITS	
1,2-Dichloroethane-d4 (Surrogate)	70-121%	80
Toluene-d8 (Surrogate)	84-138%	90
4-Bromofluorobenzene (Surrogate)	59-113%	86

J = Estimated Value Detected Below MDL

EPA SAMPLE NO.
1645.1 9/21

1E
VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

Lab Name: Princeton Testing Lab US_ARMY,FORT_MONMOUTH.

Lab Code: PTL Case No.:4760-001 SAS No.:XXX SDG No.:XXX

Matrix: (Soil/Water) Soil Lab Sample ID: 01

Sample wt/vol: 5 (g/mL) g Lab File ID: C8694

Level: (low/med) med Date Received: 09/23/94

%Moisture: not dec. 19 Date Analyzed: 09/29/94

GC Column: VOCOL ID: 0.53 mm Dilution Factor: 100

Soil Extract Vol: ul Soil Aliquot Vol: ul

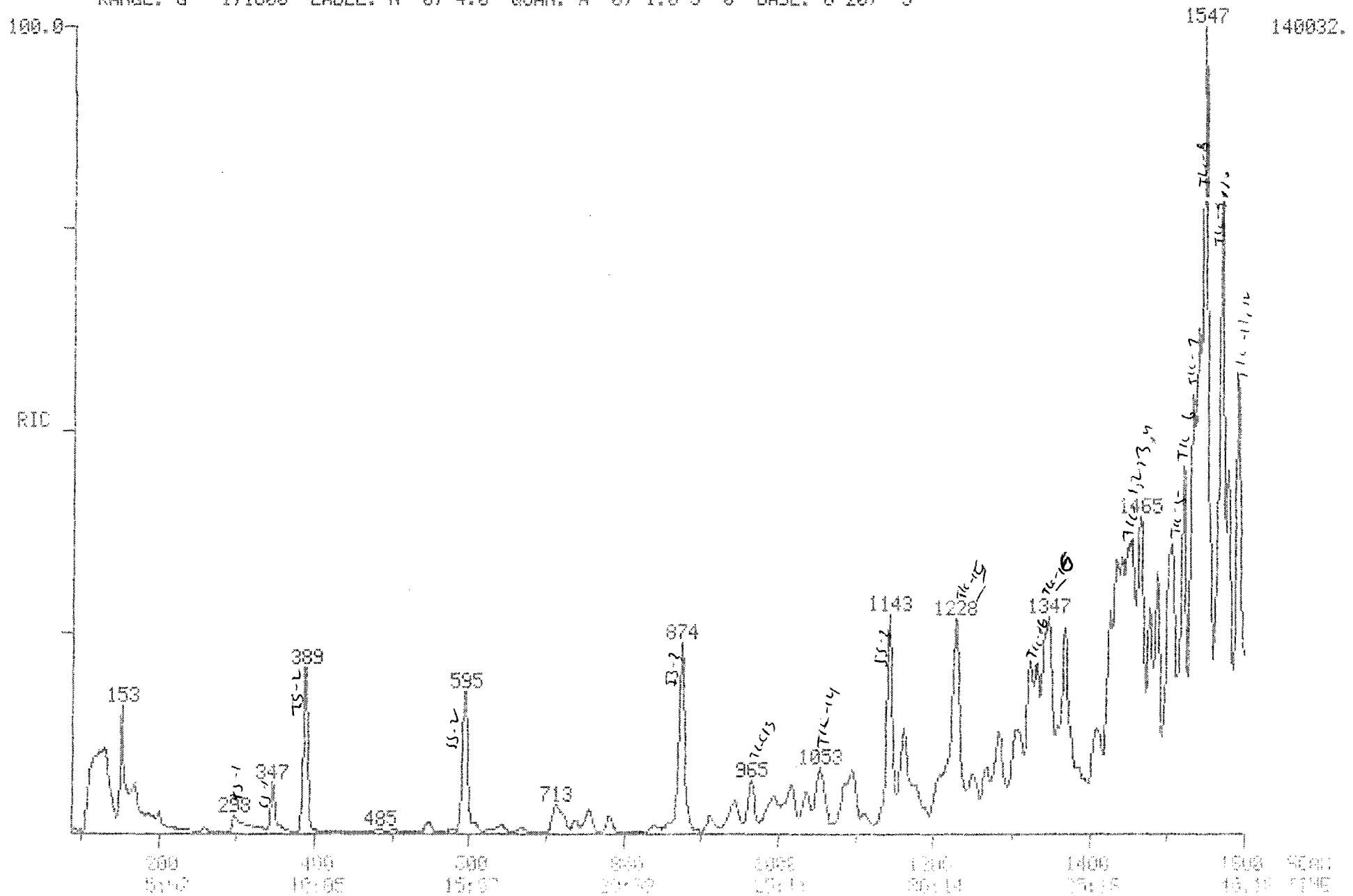
Number TICs found: 16 CONCENTRATION UNITS:
(ug/L or ug/Kg) ug/Kg

#S	CAS NUMB	COMPOUND NAME	RT	EST. CONC.	SCAN
1	0-00-0	UNKNOWN	36:30	4200	1449
2	1758-88-9	BENZENE,2-ETHYL-1,4-DIME	36:39	5300	1455
3	17301-325	UNDECANE,4,7-DIMETHYL-	36:53	9200	1464
4	16519-689	CYCLOHEXANONE,2,6-DIETHYL	36:56	6500	1466
5	1758-88-9	BENZENE,2-ETHYL-1,4-DIME	37:53	9900	1504
6	0-00-0	UNKNOWN_HYDROCARBON	38:17	6400	1520
7	2958-76-1	NAPHTHALENE,DECAHYDRO-2ME	38:34	9300	1531
8	62108-230	DECANE,2,5,6-TRIMETHYL-	38:45	16000	1538
9	0-00-0	UNKNOWN	38:58	19000	1547
10	0-00-0	UNKNOWN	39:30	14000	1568
11	527-53-7	BENZENE,1,2,3,5-TETRAMETHYL-	39:45	5300	1578
12	1758-88-9	BENZENE,2-ETHYL-1,4-DIME	40:03	8600	1590
13	0-00-0	UNKNOWN_HYDROCARBON	24:18	2800	965
14	0-00-0	UNKNOWN_HYDROCARBON	26:33	4800	1054
15	0-00-0	UNKNOWN_HYDROCARBON	30:56	16000	1228
16	0-00-0	UNKNOWN_HYDROCARBON	33:56	11000	1347

RIC
09/23/94 17:46:00
SAMPLE: 4760-001-01 1645.1 09/21/94
CONDS.: EPA METHOD 8240
RANGE: G 1,1600 LABEL: N 0, 4.0 QUAN: A 0, 1.0 J 0 BASE: U 20, 3

DATA: C8694 #1
CALI: C8694 #3

SCANS 88 TO 1600



Quantitation Report File: C8694

Data: C8694.TI
9/29/94 17:46:00
Sample: 4760-001-01 1645 1 09/21/94
Comps.: EPA METHOD 8240
Formula: 5G/10ML:100ULX100DILInstrument: FINN Weight: 0.000
Submitted by: USARMY Analyst: UC Acct. No.: 4760-001

AMOUNT=AREA * REF AMNT/(REF AREA * RESP FACT)
Resp. fac. from Library Entry

NO	Name
1	CI01 BROMOCHLOROMETHANE **INT. STD. **
2	CI10 1,4-DIFLUOROBENZENE **INT. STD. **
3	CI20 CHLOROBENZENE-D5 **INT. STD. **
4	CS15 1,2-DICHLOROETHANE-D4 **S. STD. **
5	CS05 TOLUENE-D8 **S. STD. **
6	CS10 4-BROMOFLUOROBENZENE **S. STD. **
7	C010 CHLOROMETHANE **
8	C015 BROMOMETHANE
9	C020 VINYL CHLORIDE *
10	C025 CHLOROETHANE
11	C030 METHYLENE CHLORIDE
12	C251 ACROLIN
13	C035 ACETONE
14	C252 ACRYLONITRILE
15	C040 CARBON DISULFIDE
16	C045 1,1-DICHLOROETHENE *
17	C050 1,1-DICHLOROETHANE **
18	C055 TRANS-1,2-DICHLOROETHENE
19	C000 TRICHLOROFLUOROMETHANE
20	C060 CHLOROFORM *
21	C065 1,2-DICHLOROETHANE
22	C110 2-BUTANONE
23	C115 1,1,1-TRICHLOROETHANE
24	C120 CARBON TETRACHLORIDE
25	C125 VINYL ACETATE
26	C130 BROMO DICHLOROMETHANE
27	C140 1,2-DICHLOROPROPANE *
28	C145 TRANS-1,3 DICHLOROPROPENE
29	C150 TRICHLOROETHENE
30	C155 DIBROMOCHLOROMETHANE
31	C160 1,1,2-TRICHLOROETHANE
32	C165 BENZENE
33	C143 CIS-1,3-DICHLOROPROPENE
34	C175 2-CHLOROETHYL VINYL ETHER
35	C180 BROMOFORM **
36	C220 TETRACHLOROETHENE
37	C210 2-HEXANONE
38	C205 4-METHYL 2-PENTANONE
39	C225 1,1,2,2-TETRACHLOROETHANE **
40	C230 TOLUENE *
41	C235 CHLOROBENZENE **
42	C240 ETHYL BENZENE *
43	C245 STYRENE
44	C250 M+P-XYLENES
45	C253 1,3-DICHLOROBENZENE
46	C254 1,4-DICHLOROBENZENE
47	C255 1,2-DICHLOROBENZENE

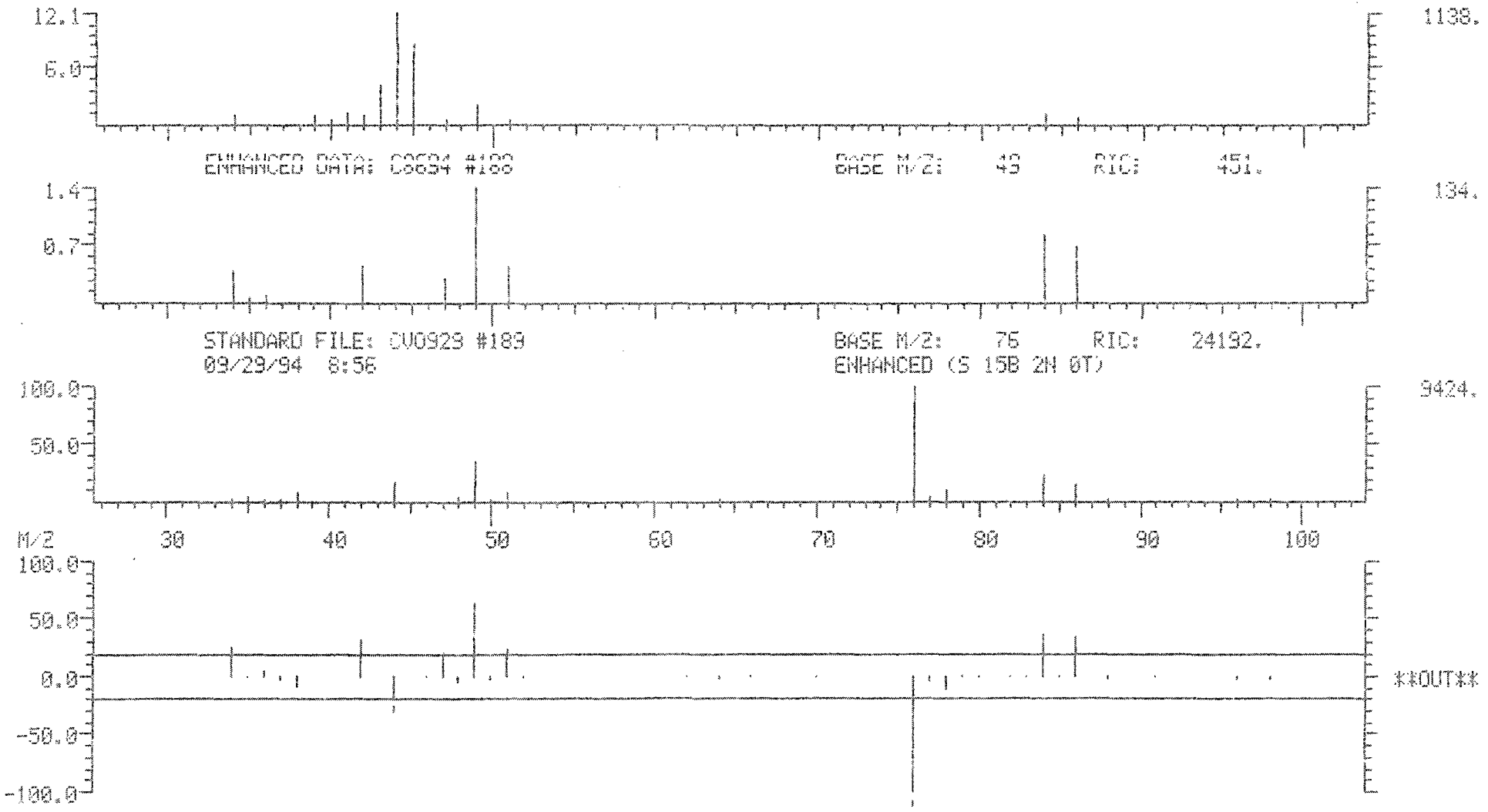
1864

No	Name
48	C250 O-XYLENE
49	BENZENE-D6 **S. STD. **

No	m/z	Scan	Time	Ref	RRT	Meth	Area(Hght)	Amount	%Tot
1	49	298	7:30	1	1.000	M XX	10152.	50.000 NG	17.00
2	114	388	9:46	2	1.000	A BB	68901.	50.000 NG	17.00
3	117	874	22:01	3	1.000	A BB	56616.	50.000 NG	17.00
4	65	347	8:44	1	1.164	A BB	13805.	40.204 NG	13.67
5	98	595	14:59	3	0.681	A BB	67904.	44.753 NG	15.22
6	95	1143	28:48	3	1.308	A BB	47236.	43.084 NG	14.65
7	NOT FOUND								
8	NOT FOUND								
9	NOT FOUND								
10	NOT FOUND								
11	49	190	4:47	1	0.638	A BV	1830.	3.692 NG	1.26
12	NOT FOUND								
13	NOT FOUND								
14	NOT FOUND								
15	NOT FOUND								
16	NOT FOUND								
17	NOT FOUND								
18	NOT FOUND								
19	NOT FOUND								
20	NOT FOUND								
21	NOT FOUND								
22	43	259	6:31	2	0.668	A BB	2487.	11.804 NG	4.01
23	NOT FOUND								
24	NOT FOUND								
25	43	225	5:40	2	0.580	A BV	121.	0.126 NG	0.04
26	NOT FOUND								
27	NOT FOUND								
28	NOT FOUND								
29	NOT FOUND								
30	NOT FOUND								
31	NOT FOUND								
32	NOT FOUND								
33	NOT FOUND								
34	NOT FOUND								
35	NOT FOUND								
36	NOT FOUND								
37	NOT FOUND								
38	NOT FOUND								
39	NOT FOUND								
40	91	607	15:17	3	0.695	A BB	1107.	0.631 NG	0.21
41	NOT FOUND								
42	NOT FOUND								
43	NOT FOUND								
44	NOT FOUND								
45	NOT FOUND								
46	NOT FOUND								
47	NOT FOUND								
48	NOT FOUND								
49	NOT FOUND								

DATA FILE: C8694 #188 STANDARD FILE: CV0929 #189
TARGET COMPOUND COMPARISON CALI: C8694 #3
COMPOUND: C030 METHYLENE CHLORIDE

RAW DATA: C8694 #188 BASE M/Z: 44 RIC: 3276.
09/29/94 17:46

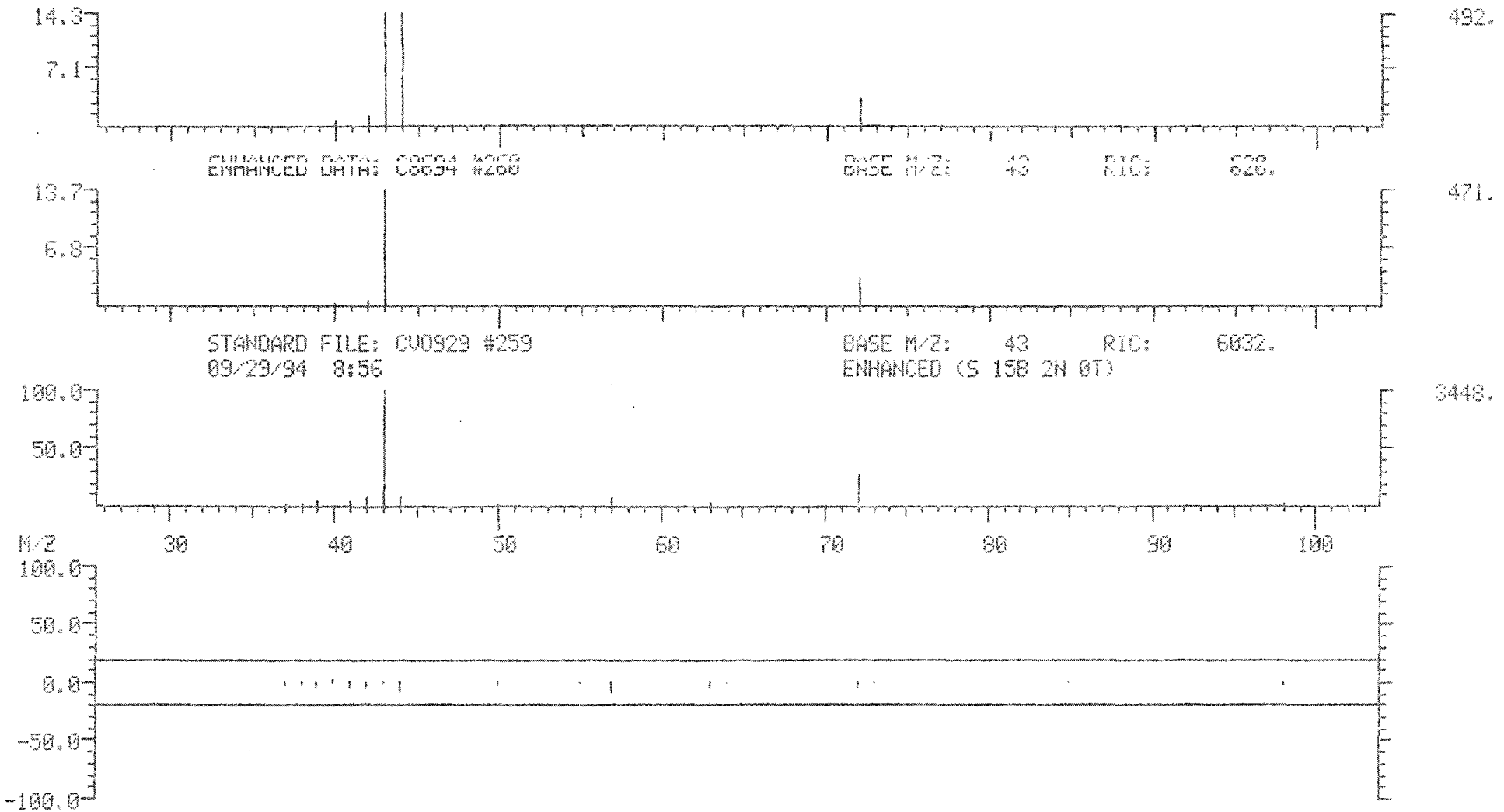


DATA FILE: C8694 #260
TARGET COMPOUND COMPARISON
COMPOUND: C110 2-BUTANONE

STANDARD FILE: CV0929 #259
CALI: C8694 #3

RAW DATA: C8694 #260
09/29/94 17:45

BASE M/Z: 43 RIC: 1154.



PROCEDURE: FILTER/TIC DIAGNOSTIC REPORT 9/29/94 18:53:10

DATA FILE: C8694

USER SCAN PARAMETERS	
MAX NUMBER TICS:	15
LIBRARY TABLE ENTRIES:	539
SCAN TOLERANCE :	2
MIN RIC HT. (%):	10
START SCAN :	1
END SCAN :	1600
TIC THRESHOLD :	600

METHOD LIBRARY & LISTS	
TIC I.S. LIBRARY:	LIBRARYLS
NBS SEARCH PROC :	SERLIB
PEAK FINDER PROC:	VOME
TCA I.S. LL :	LS
FILE NAME LIST :	TCAREF2

TARGET COMPOUND ANALYSIS:

TARGETS (QUAN LIST)	IS PEAKS	TOTAL TARGET PEAKS
7	3	10

FILTER PROCESSING:

<----------REJECT PEAKS---------->							
TOTAL PEAKS	< 1ST SCAN	> LAST SCAN	< MIN RIC HT	< SCAN TOL	> MAX # PEAKS	TOTAL REJECTS	TOTAL TICS
46	0	0	0	6	25	31	15

TIC PROCESSING:

#	SCAN#	PURITY	FIT	MW	COMPOUND NAME [BEFORE TIC THRESHOLD]
1	82	431	453	42	BORON, CARBONYLTRIHYDRO-, (T-4
2	84	994	1000	44	CARBON DIOXIDE (ACN)
3	1449	495	861	168	1-NONENE, 4,6,8-TRIMETHYL-
4	1454	732	865	134	BENZENE, 2-ETHYL-1,4-DIMETHYL-
5	1463	823	917	184	UNDECANE, 4,7-DIMETHYL-
6	1467	662	810	154	CYCLOHEXANONE, 2,6-DIETHYL-
7	1506	838	968	134	BENZENE, 2-ETHYL-1,4-DIMETHYL-
8	1520	846	990	134	BENZENE, 2-ETHYL-1,4-DIMETHYL-
9	1531	608	936	152	NAPHTHALENE, DECAHYDRO-2-METHY
10	1533	432	891	152	NAPHTHALENE, DECAHYDRO-2-METHY
11	1539	649	910	184	DECANE, 2,5,6-TRIMETHYL-
12	1547	511	860	256	1-HEPTADECANOL
13	1568	495	902	240	OXIRANE, TETRADECYL-
14	1578	712	958	134	BENZENE, 1,2,3,5-TETRAMETHYL-
15	1590	842	975	134	BENZENE, 2-ETHYL-1,4-DIMETHYL-

Identification Report File C8694

Data: C8694.TI
09/29/94 17:46:00
Sample: 4760-001-01 1645.1 09/21/94
Comm.: EPA METHOD 8240
Formula: 5G/10ML:100ULX100DILInstrument: FINN Weight: 0.000
Submitted by: USARMY Analyst: UC Acct. No.: 4760-001

AMOUNT=AREA * REF AMNT/(REF AREA * RESP FACT)
Resp. fac. from Library Entry

No	CAS #	Name
1	0-00-0	C101 BROMOCHLOROMETHANE **INT.STD.**
2	0-00-0	C110 1,4-DIFLUOROBENZENE **INT.STD.**
3	0-00-0	C120 CHLOROBENZENE-D5 **INT.STD.**
4	0-00-0	UNKNOWN
5	124-38-9	CARBON DIOXIDE (ACN)
6	0-00-0	UNKNOWN
7	1758-88-9	BENZENE, 2-ETHYL-1,4-DIMETHYL-
8	17301-32-5	UNDECANE, 4,7-DIMETHYL-
9	16519-68-9	CYCLOHEXANONE, 2,6-DIETHYL-
10	1758-88-9	BENZENE, 2-ETHYL-1,4-DIMETHYL-
11	1758-88-9	BENZENE, 2-ETHYL-1,4-DIMETHYL-
12	2958-76-1	NAPHTHALENE, DECAHYDRO-2-METHYL-
13	0-00-0	UNKNOWN
14	62108-23-0	DECANE, 2,5,6-TRIMETHYL-
15	0-00-0	UNKNOWN
16	0-00-0	UNKNOWN
17	527-53-7	BENZENE, 1,2,3,5-TETRAMETHYL-
18	1758-88-9	BENZENE, 2-ETHYL-1,4-DIMETHYL-

No	m/z	Scan	Time	Ref	RRT	Meth	Area(Hght)	Amount	%Tot
1	TOT	298	7:30	0	ISINV	A BB	27554.	****** UG/L	00.00
2	TOT	388	9:46	0	ISINV	A BB	152265.	****** UG/L	00.00
3	TOT	874	22:01	0	ISINV	A BB	147116.	****** UG/L	00.00
4	TOT	82	2:04	1	0.275	A BB	368441.	668.568	36.95
5	TOT	84	2:07	1	0.282	A BB	83163.	150.908	8.34
6	TOT	1449	36:30	3	1.658	A VB	99170.	33.705	1.86
7	TOT	1455	36:39	3	1.665	A VB	126607.	43.030	2.38
8	TOT	1464	36:53	3	1.675	A BV	218076.	74.117	4.10
9	TOT	1466	36:56	3	1.677	A BV	155584.	52.878	2.92
10	TOT	1504	37:53	3	1.721	A BB	235829.	80.151	4.43
11	TOT	1520	38:17	3	1.739	A BB	152142.	51.708	2.86
12	TOT	1531	38:34	3	1.752	A BV	221133.	75.156	4.15
13	TOT	1531	38:34	3	1.752	A BV	221133.	75.156	4.15
14	TOT	1538	38:45	3	1.760	A BV	377421.	128.274	7.09
15	TOT	1547	38:58	3	1.770	A VB	450144.	152.990	8.46
16	TOT	1568	39:30	3	1.794	A BV	322966.	109.766	6.07
17	TOT	1578	39:45	3	1.805	A VB	127205.	43.233	2.39
18	TOT	1590	40:03	3	1.819	A BB	205032.	69.684	3.85

No	Ret(L)	Ratio	RRT(L)	Ratio	Amnt	Amnt(L)	R.Fac	R.Fac(L)	Ratio
1	9:13	0.81	1.000						
2	19:21	0.51	1.000						
3	23:54	0.92	1.000						

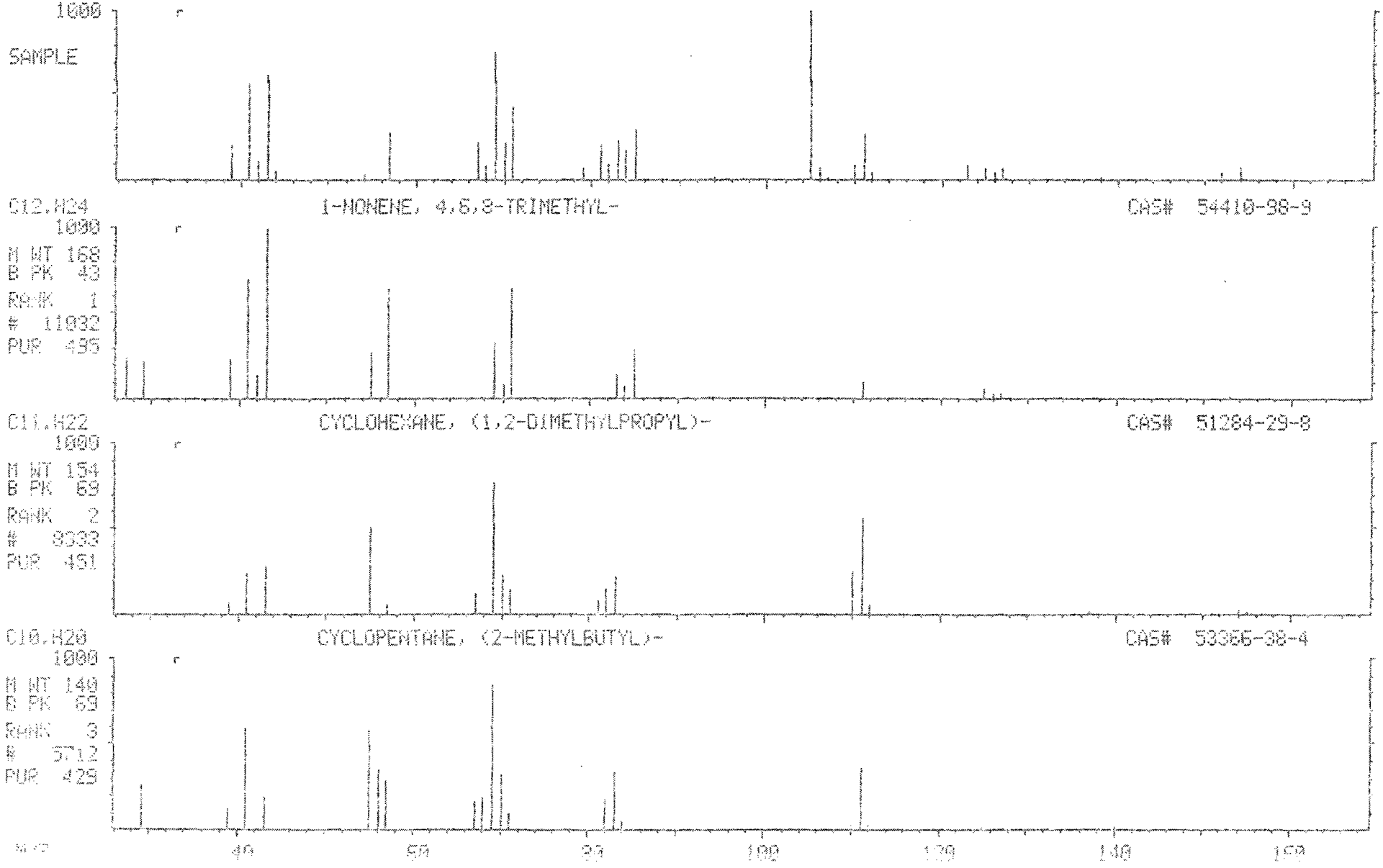
C0647

Mo	Ret(L) Ratio	RRT(L) Ratio	Amnt	Amnt(L)	R.Fac	R.Fac(L)	Ratio
4			668.57	1.00	668.570	1.000	668.57
5			150.91	1.00	150.908	1.000	150.91
6			33.70	1.00	33.705	1.000	33.70
7			43.03	1.00	43.030	1.000	43.03
8			74.12	1.00	74.117	1.000	74.12
9			52.88	1.00	52.878	1.000	52.88
10			80.15	1.00	80.151	1.000	80.15
11			51.71	1.00	51.708	1.000	51.71
12			75.16	1.00	75.156	1.000	75.16
13			75.16	1.00	75.156	1.000	75.16
14			128.27	1.00	128.274	1.000	128.27
15			152.99	1.00	152.990	1.000	152.99
16			109.77	1.00	109.766	1.000	109.77
17			43.23	1.00	43.233	1.000	43.23
18			69.68	1.00	69.684	1.000	69.68

MID LIBRARY SEARCH (LIBRARYNB)
09/29/94 17:46:00 + 36:30
SAMPLE: 4760-001-01 1645.1 09/21/94
CONDS.: EPA METHOD 8240
ENHANCED (S 15B 2N 0T)

DATA: C8694 #1449
CALI: C8694 # 3

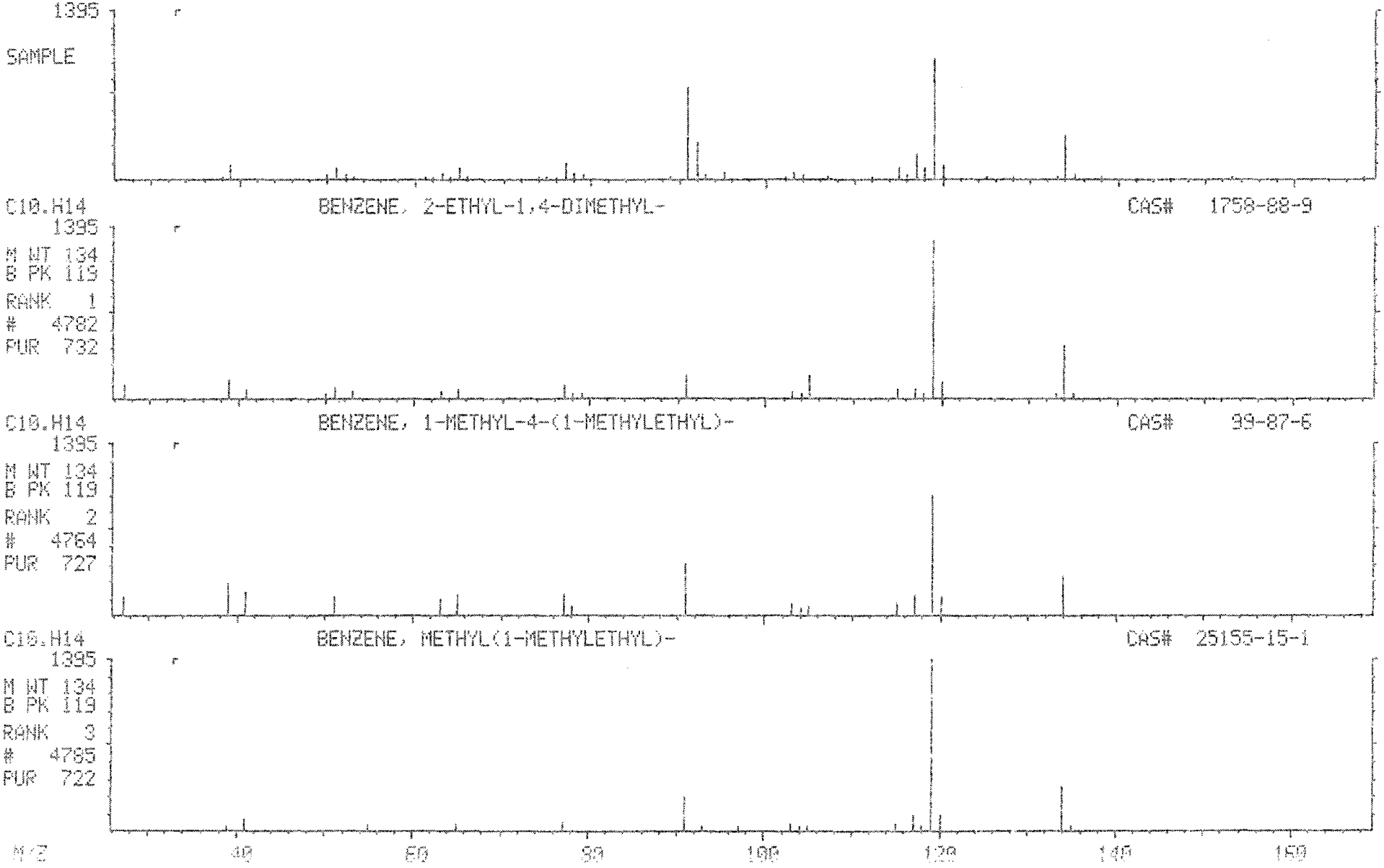
BASE M/Z: 105
RIC: 15712.



MID LIBRARY SEARCH (LIBRARYNB)
09/29/94 17:45:00 + 36:38
SAMPLE: 4760-001-01 1645.1 09/21/94
CONDS.: EPA METHOD 8240
ENHANCED (S 15B 2N 0T)

DATA: C8694 #1454
CALI: C8694 # 3

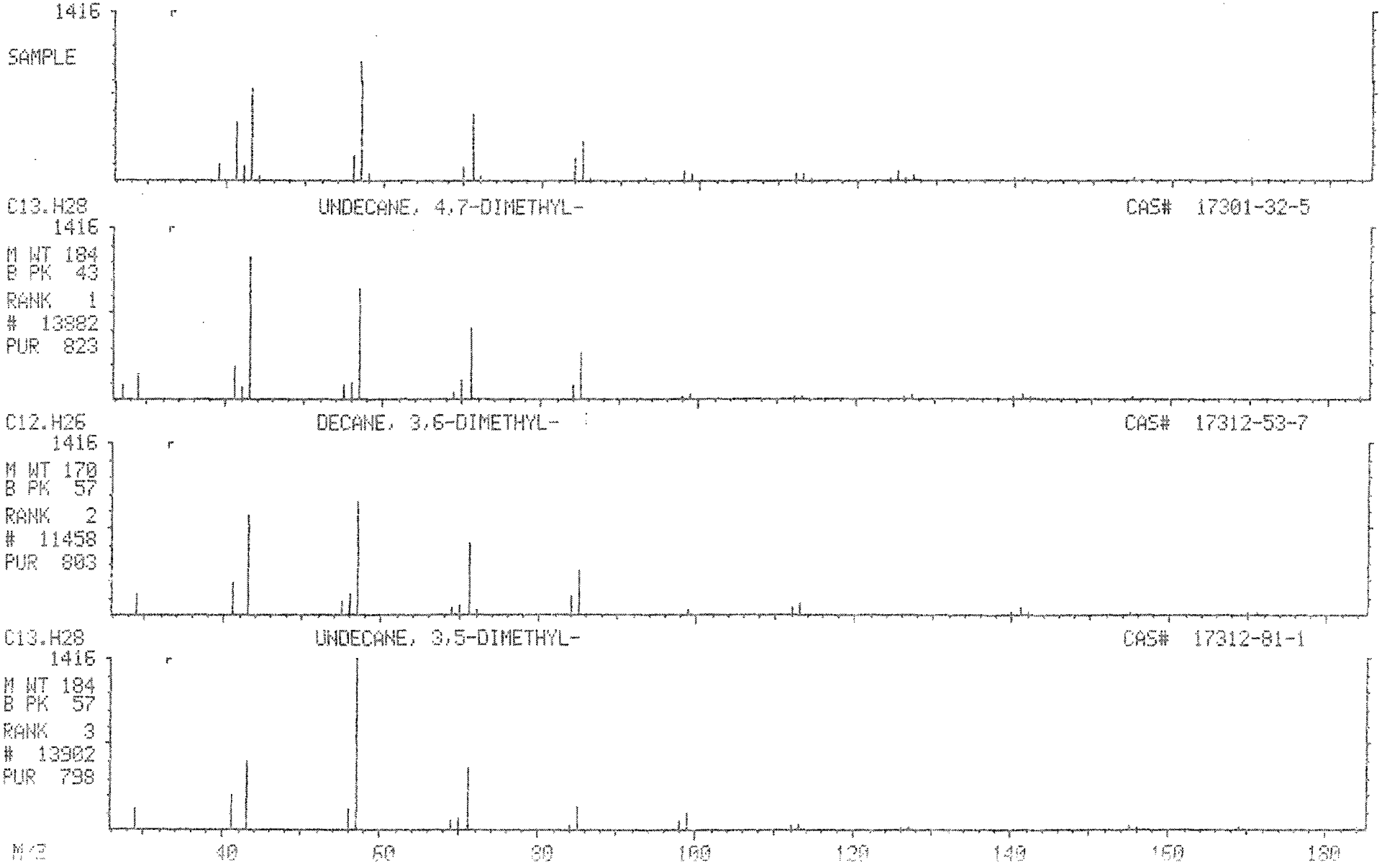
BASE M/Z: 119
RIC: 19584.



MID LIBRARY SEARCH (LIBRARYNB)
09/29/94 17:46:00 + 36:51
SAMPLE: 4760-001-01 1645.1 09/21/94
CONDS.: EPA METHOD 8240
ENHANCED (S 15B 2N 0T)

DATA: C8694 #1463
CALI: C8694 # 3

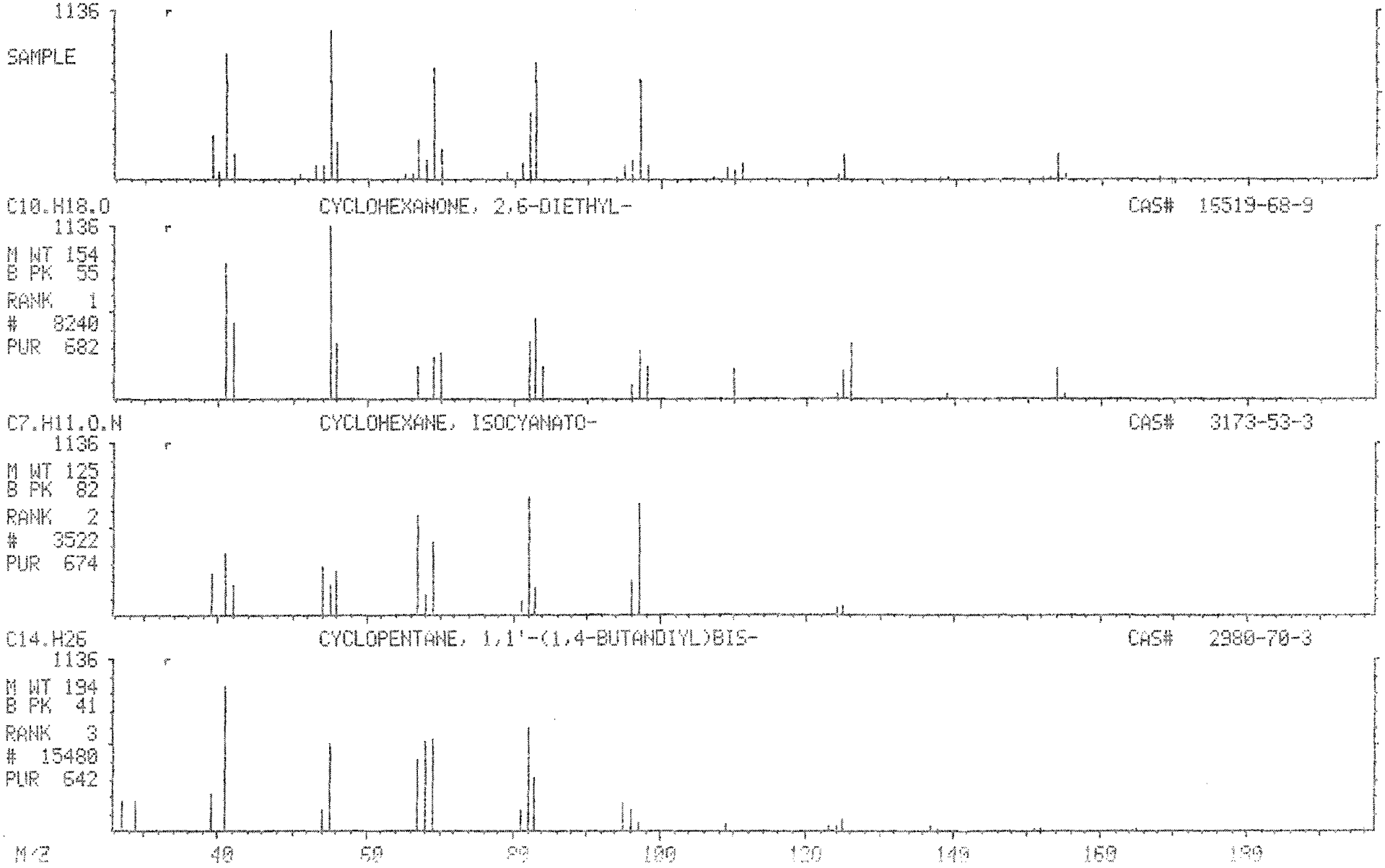
BASE M/Z: 57
RIC: 22848.



MID LIBRARY SEARCH (LIBRARYNB)
09/29/94 17:46:00 + 36:57
SAMPLE: 4760-001-01 1645.1 09/21/94
CONDS.: EPA METHOD 8240
ENHANCED (S 15B 2N 0T)

DATA: C8694 #1467
CALI: C8694 # 3

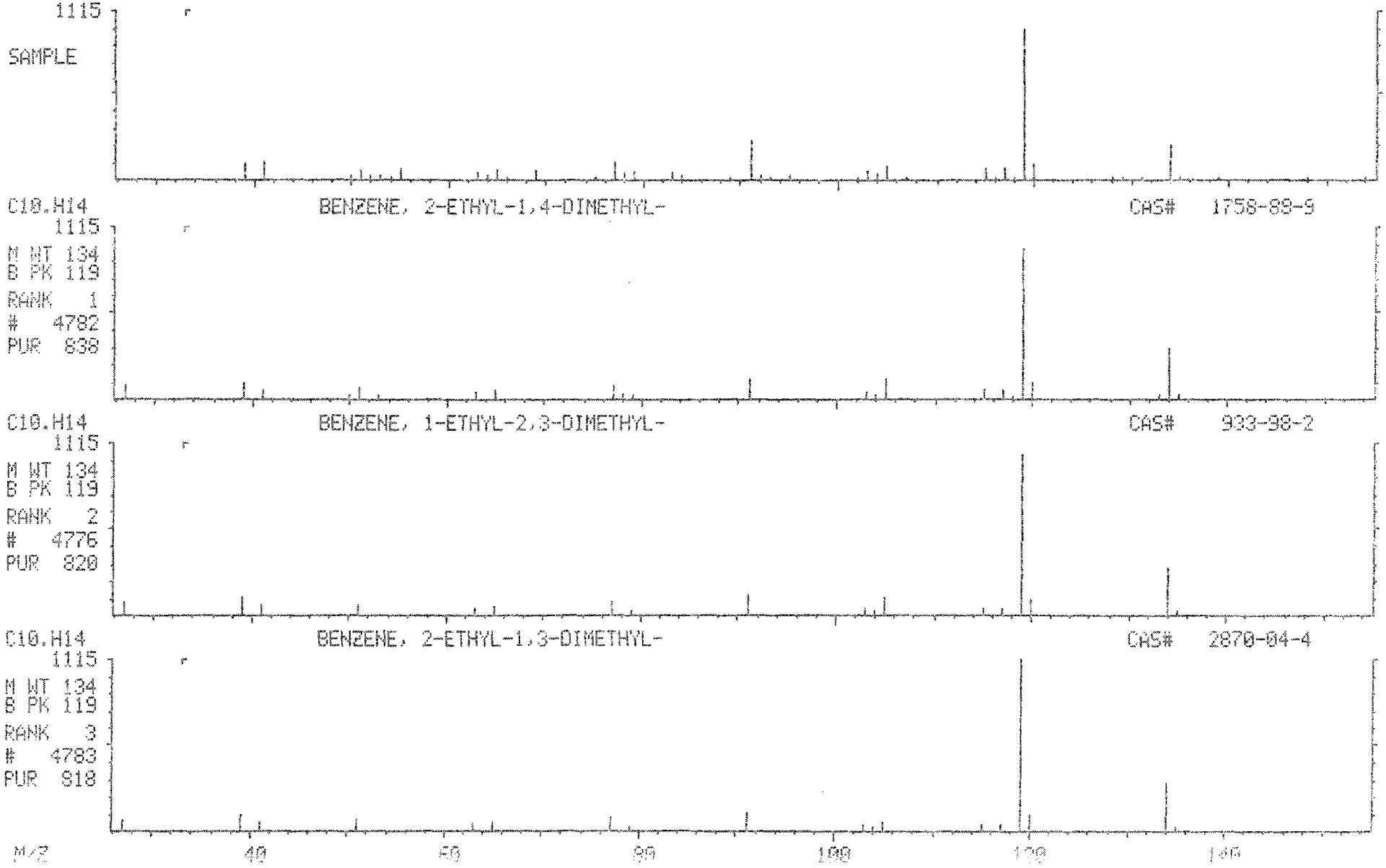
BASE M/Z: 55
RIC: 18784.



MID LIBRARY SEARCH (LIBRARYNB)
09/29/94 17:46:00 + 37:56
SAMPLE: 4760-001-01 1645.1 09/21/94
CONDS.: EPA METHOD 8240
ENHANCED (S 15B 2N 0T)

DATA: C8694 #1506
CALI: C8694 # 3

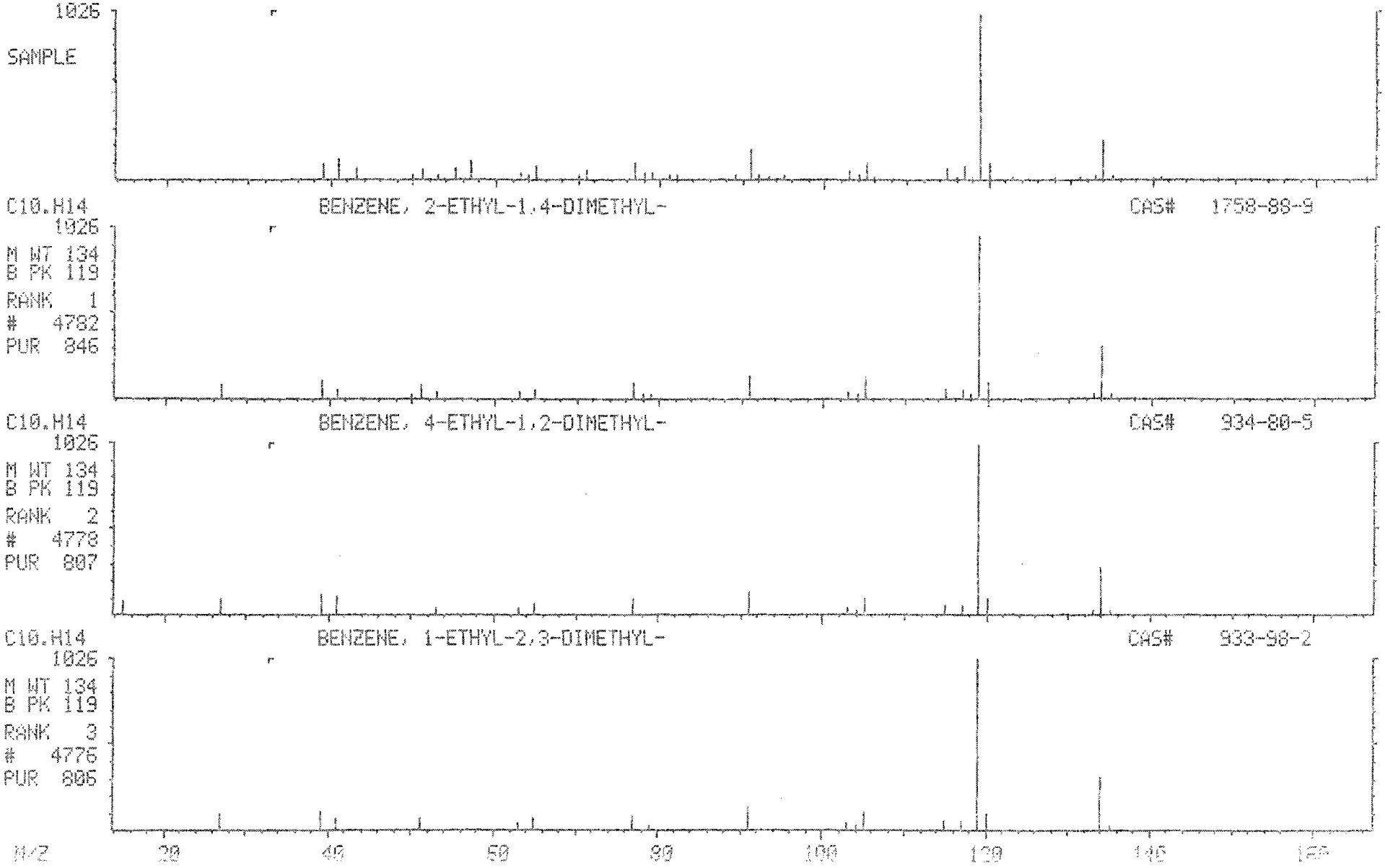
BASE M/Z: 119
RIC: 19648.



MID LIBRARY SEARCH (LIBRARYNB)
09/29/94 17:46:00 + 38:17
SAMPLE: 4760-001-01 1645.1 09/21/94
CONDS.: EPA METHOD 8240
ENHANCED (S 15B 2N 0T)

DATA: C8694 #1520
CALI: C8694 # 3

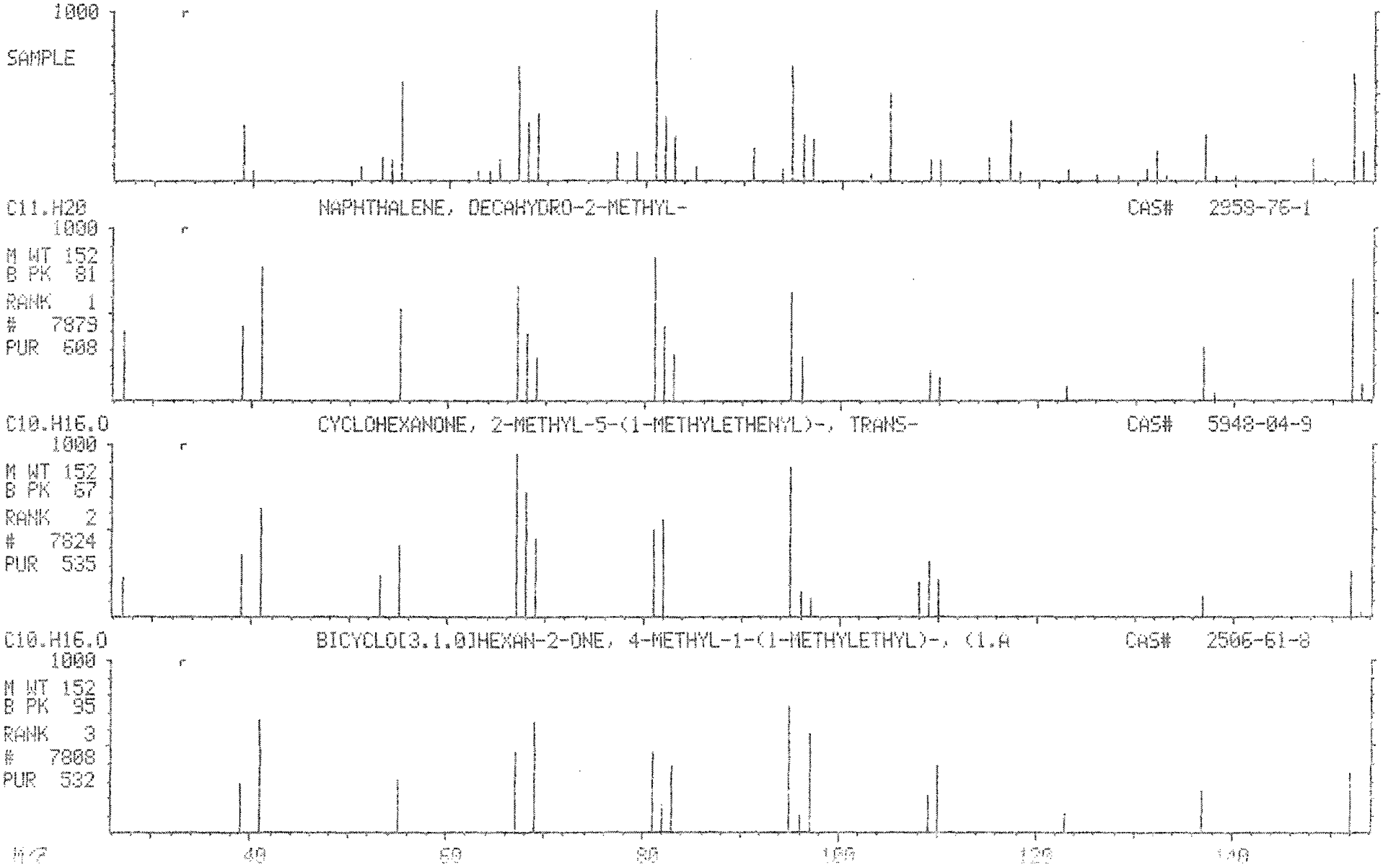
BASE M/Z: 119
RIC: 35392.



MID LIBRARY SEARCH (LIBRARYNB)
09/29/94 17:46:00 + 38:34
SAMPLE: 4760-001-01 1645.1 09/21/94
CONDS.: EPA METHOD 8240
ENHANCED (S 15B 2N 0T)

DATA: C8694 #1531
CALI: C8694 # 3

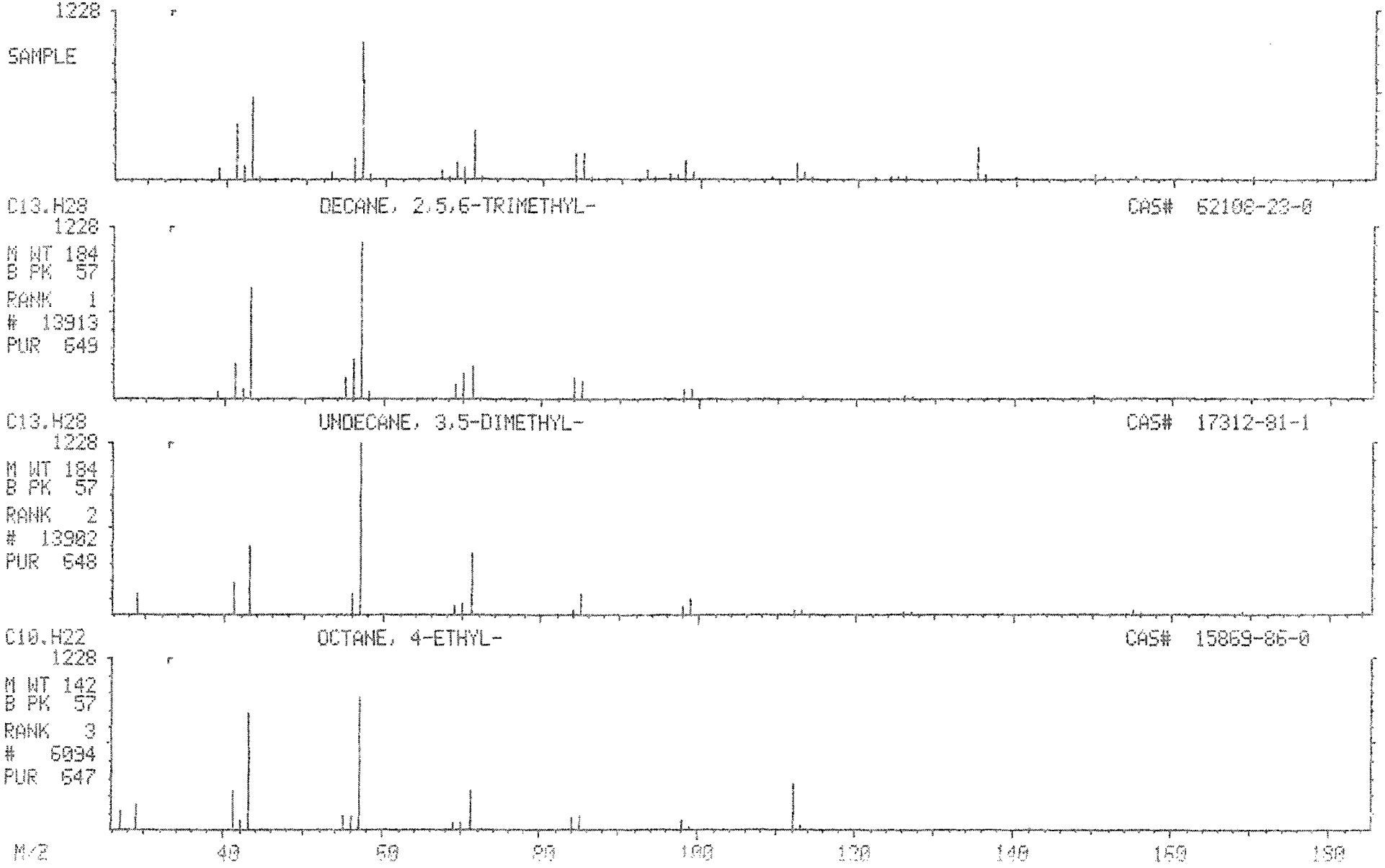
BASE M/Z: 81
RIC: 27456.



MID LIBRARY SEARCH (LIBRARYNB)
09/29/94 17:46:00 + 38:46
SAMPLE: 4760-001-01 1645.1 09/21/94
CONDS.: EPA METHOD 8240
ENHANCED (S 15B 2N 0T)

DATA: C8694 #1539
CALI: C8694 # 3

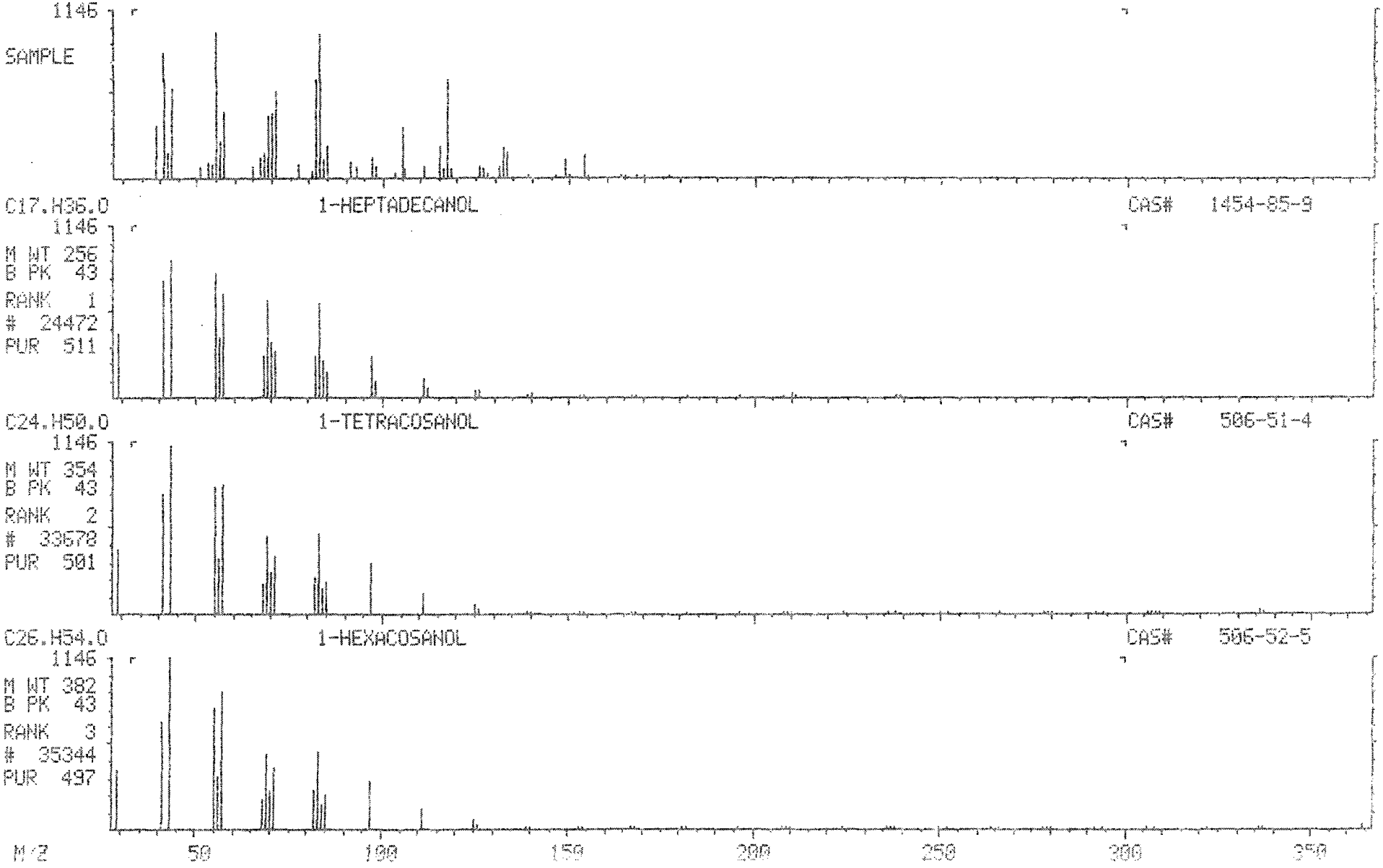
BASE M/Z: 57
RIC: 28928.



MID LIBRARY SEARCH (LIBRARYNB)
09/29/94 17:46:00 + 38:58
SAMPLE: 4760-001-01 1645.1 09/21/94
CONDS.: EPA METHOD 8240
ENHANCED (S 15B 2N 0T)

DATA: C8694 #1547
CALI: C8694 # 3

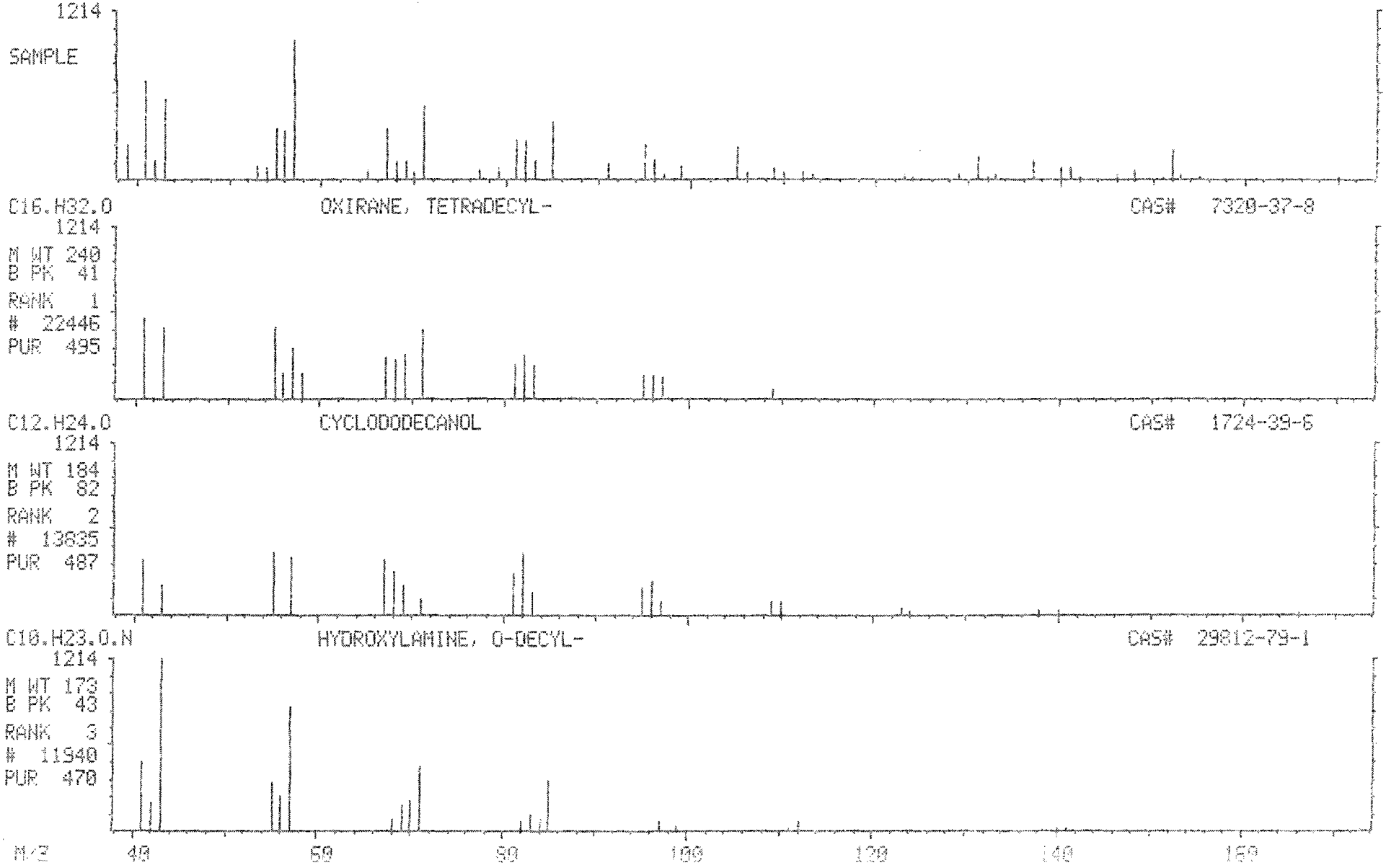
BASE M/Z: 55
RIC: 77056.



MID LIBRARY SEARCH (LIBRARYNB)
09/29/94 17:46:00 + 39:30
SAMPLE: 4760-001-01 1645.1 09/21/94
CONDS.: EPA METHOD 8240
ENHANCED (S 15B 2N 0T)

DATA: C8694 #1568
CALI: C8694 # 3

BASE M/Z: 57
RIC: 62144.

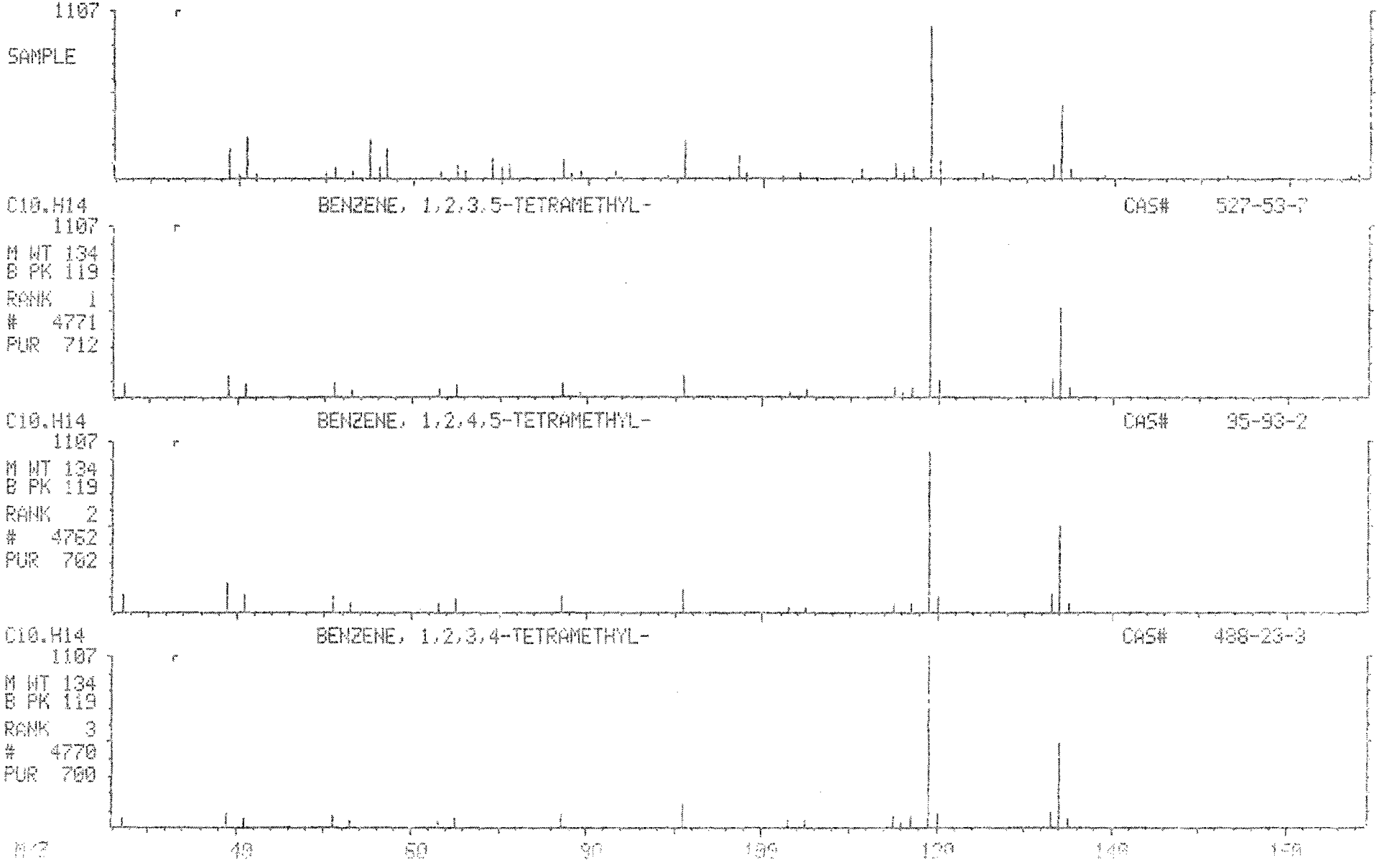


50

MID LIBRARY SEARCH (LIBRARYNB)
09/29/94 17:46:00 + 39:45
SAMPLE: 4760-001-01 1645.1 09/21/94
CONDS.: EPA METHOD 8240
ENHANCED (S 15B 2N 0T)

DATA: C8694 #1570
CALI: C8694 # 3

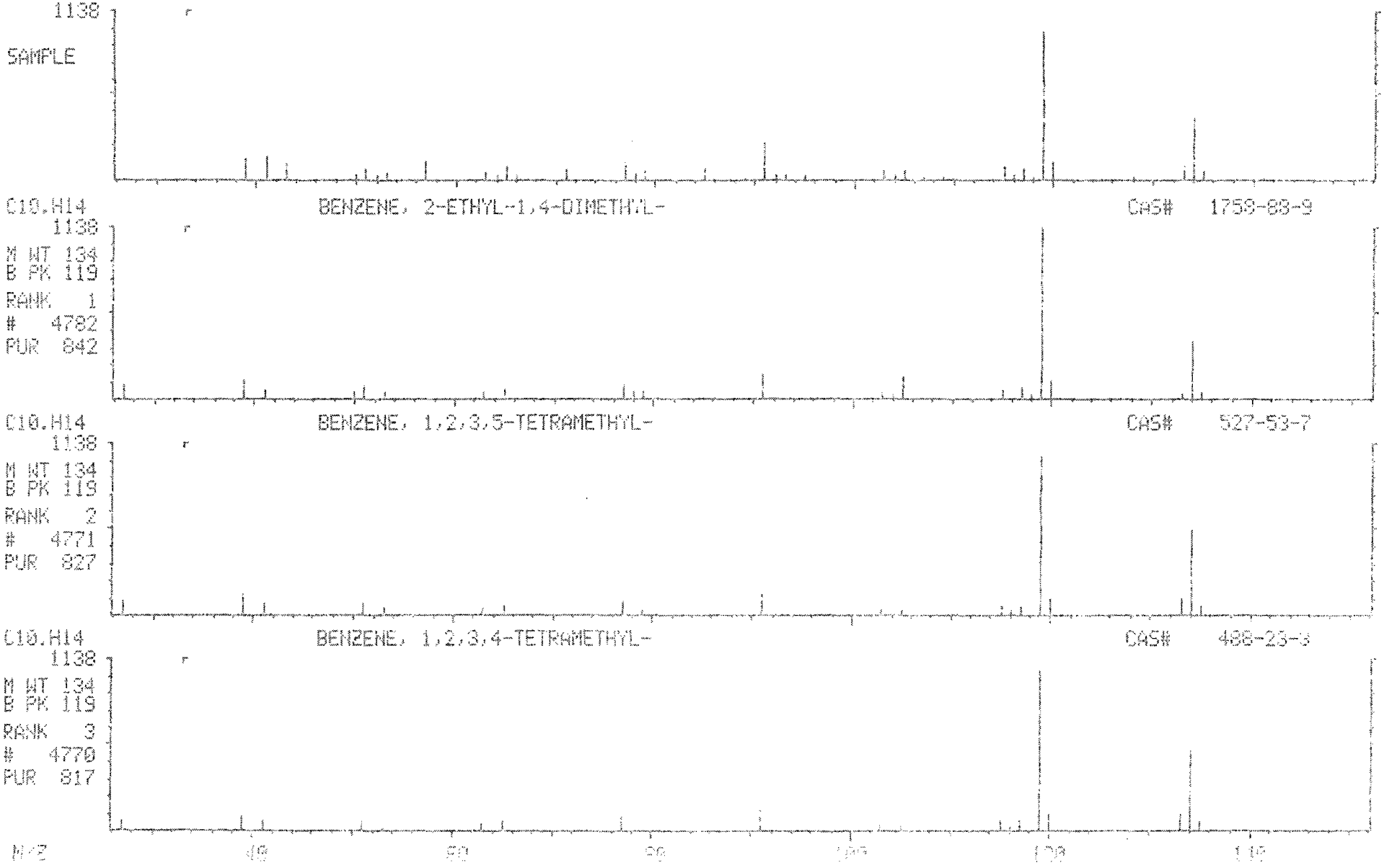
BASE M/Z: 119
RIC: 25920.



MID LIBRARY SEARCH (LIBRARYNB)
09/29/94 17:46:00 + 40:03
SAMPLE: 4760-001-01 1645.1 03/21/94
CONDS.: EPA METHOD 8240
ENHANCED (S 15B 2N 0T)

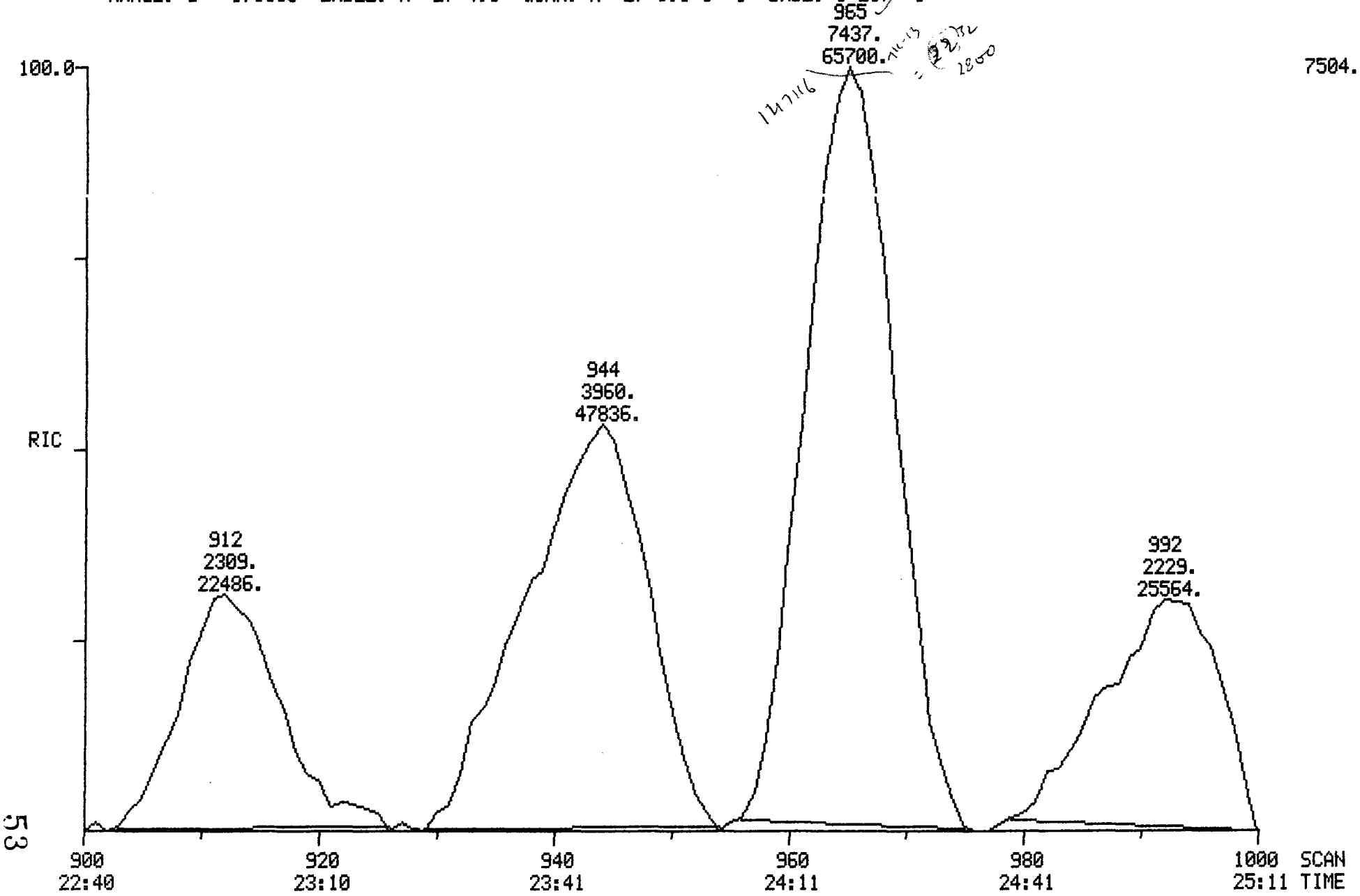
DATA: C8694 #1530
CALI: C8694 # 3

BASE M/Z: 113
RIC: 43264.



RIC
09/29/94 17:46:00
SAMPLE: 4760-001-01 1645.1 09/21/94
CONDS.: EPA METHOD 8240
RANGE: G 1,1600 LABEL: N 2, 4.0 QUAN: A 2, 1.0 J 0 BASE: U 20, 3

DATA: C8694 #1228 SCANS 900 TO 1000
CALI: C8694 #3



Library Search Data: C8694 # 965 Base m/z: 57
09/29/94 17:46:00 + 24:18 Cali: C8694 # 3 RIC: 5640.
Sample: 4760-001-01 1645.1 09/21/94
Conds.: EPA METHOD 8240
Enhanced (S 15B 2N 0T)

42223 spectra in LIBRARYNB searched for maximum FIT
171 matched at least 8 of the 16 largest peaks in the unknown

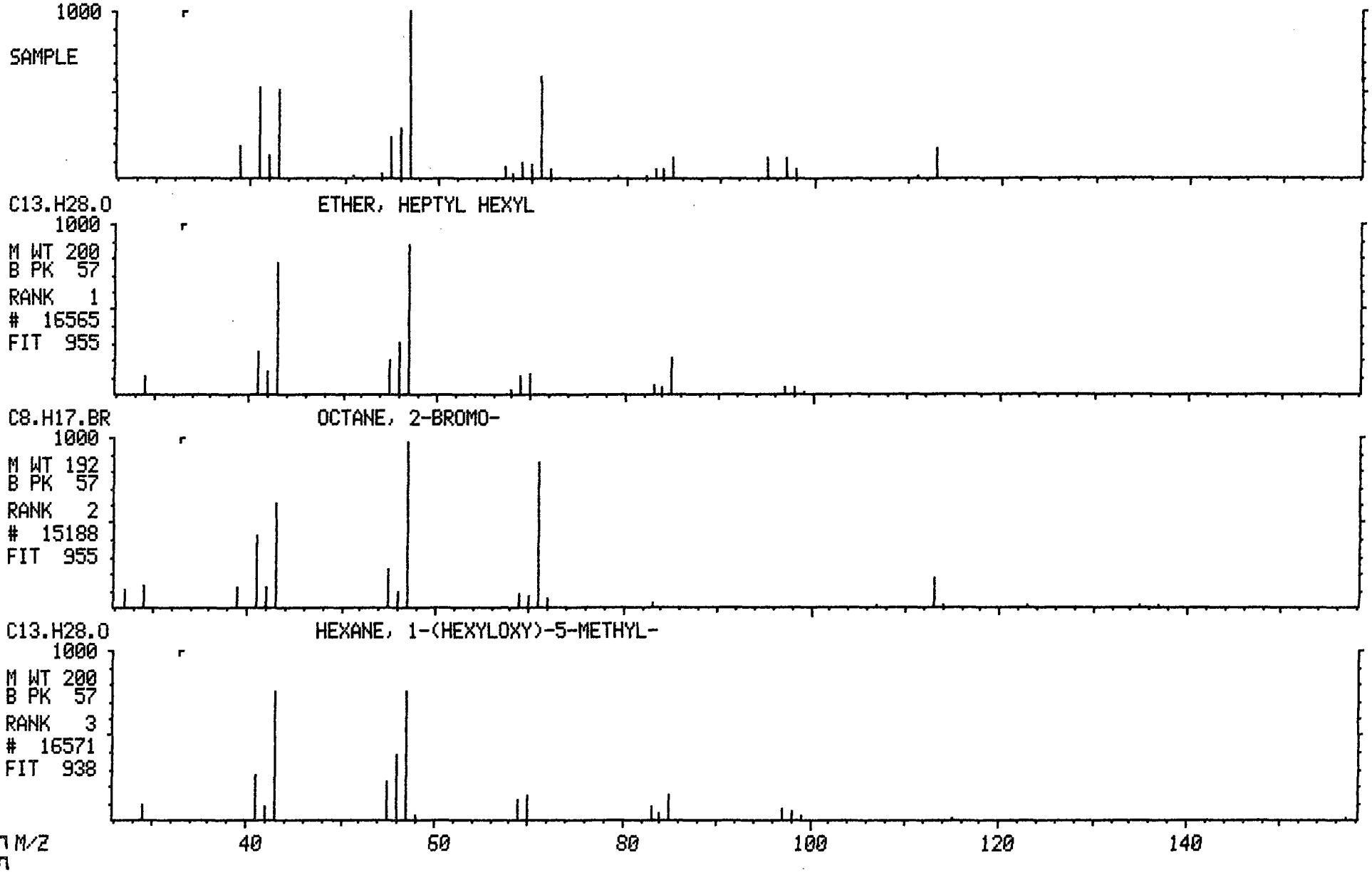
Rank	In.	Name
1	16565	ETHER, HEPTYL HEXYL
2	15188	OCTANE, 2-BROMO-
3	16571	HEXANE, 1-(HEXYLOXY)-5-METHYL-
4	6650	1-PENTANOL, 4-METHYL-2-PROPYL-
5	9069	1-HEPTANOL, 2-PROPYL-

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C13.H28.O	200	57	635	955	641
2	C8.H17.BR	192	57	775	955	791
3	C13.H28.O	200	57	619	938	638
4	C9.H20.O	144	57	780	930	783
5	C10.H22.O	158	43	755	920	773

MID LIBRARY SEARCH (LIBRARYNB)
09/29/94 17:46:00 + 24:18
SAMPLE: 4760-001-01 1645.1 09/21/94
CONDS.: EPA METHOD 8240
ENHANCED (S 15B 2N 0T)

DATA: C8694 # 965
CALI: C8694 # 3

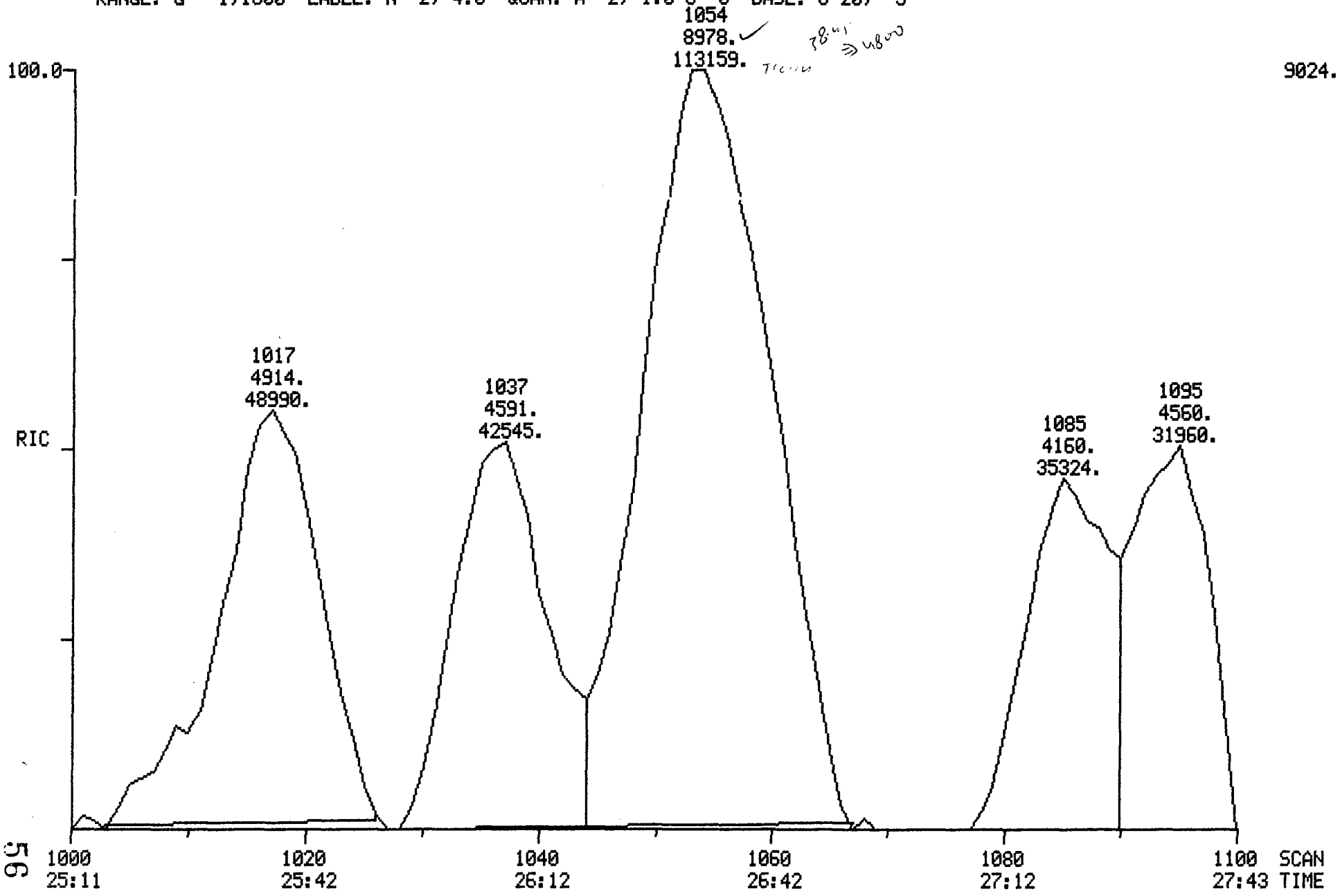
BASE M/Z: 57
RIC: 5640.



RIC
09/29/94 17:46:00
SAMPLE: 4760-001-01 1645.1 09/21/94
CONDS.: EPA METHOD 8240
RANGE: G 1,1600 LABEL: N 2, 4.0 QUAN: A 2, 1.0 J 0 BASE: U 20, 3

DATA: C8694 #1228
CALI: C8694 #3

SCANS 1000 TO 1100



Library Search Data: C8694 #1054 Base m/z: 57
09/29/94 17:46:00 + 26:33 Cali: C8694 # 3 RIC: 5120.
Sample: 4760-001-01 1645.1 09/21/94
Conds.: EPA METHOD 8240
Enhanced (S 15B 2N 0T)

42223 spectra in LIBRARYNB searched for maximum FIT
279 matched at least 7 of the 16 largest peaks in the unknown

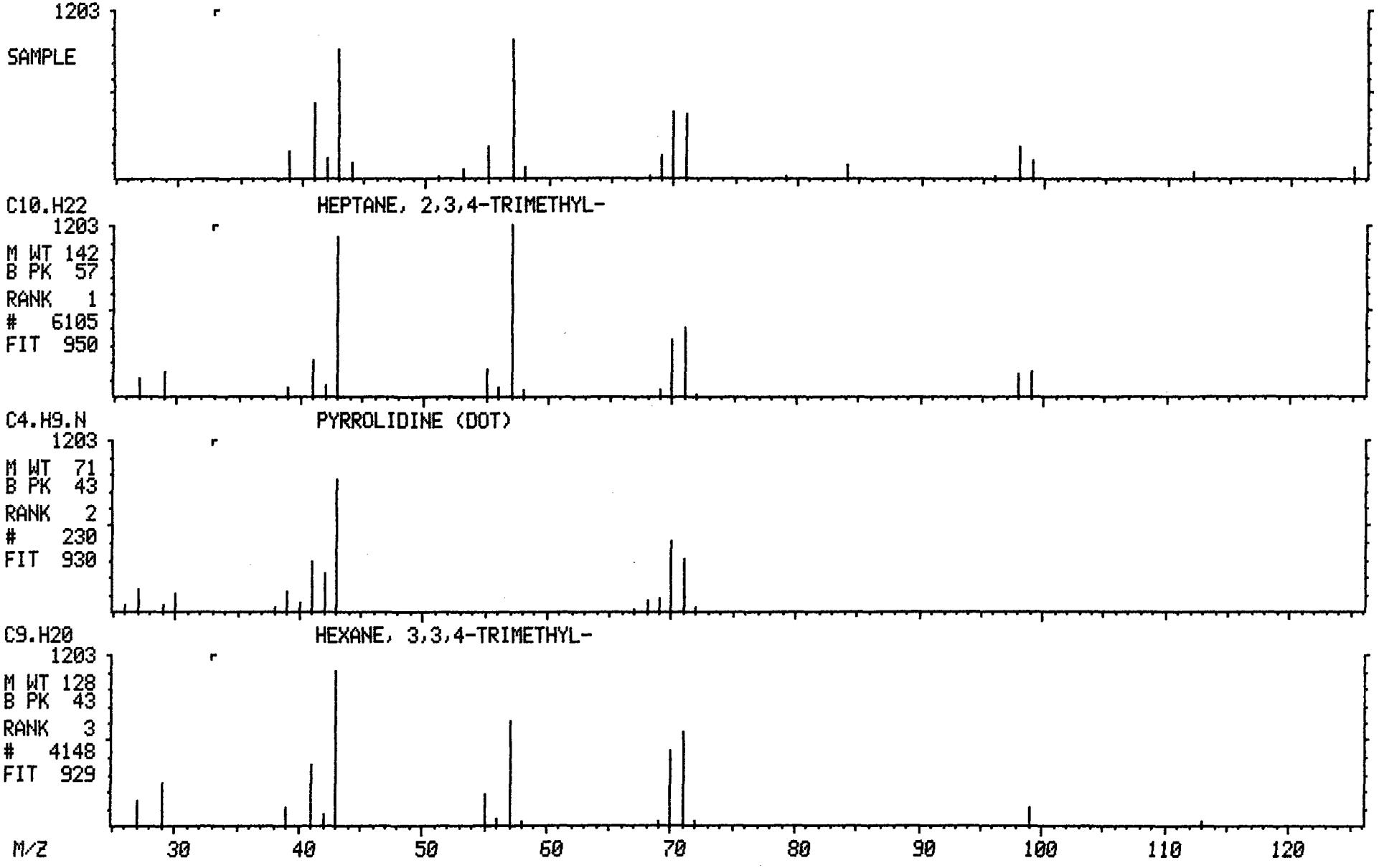
Rank	In.	Name
1	6105	HEPTANE, 2,3,4-TRIMETHYL-
2	230	PYRROLIDINE (DOT)
3	4148	HEXANE, 3,3,4-TRIMETHYL-
4	4142	PENTANE, 2,2,3,3-TETRAMETHYL-
5	4147	HEXANE, 2,4,4-TRIMETHYL-

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C10.H22	142	57	837	950	854
2	C4.H9.N	71	43	481	930	504
3	C9.H20	128	43	749	929	777
4	C9.H20	128	57	716	912	766
5	C9.H20	128	43	764	907	788

MID LIBRARY SEARCH (LIBRARYNB)
09/29/94 17:46:00 + 26:33
SAMPLE: 4760-001-01 1645.1 09/21/94
CONDS.: EPA METHOD 8240
ENHANCED (S 15B 2N 0T)

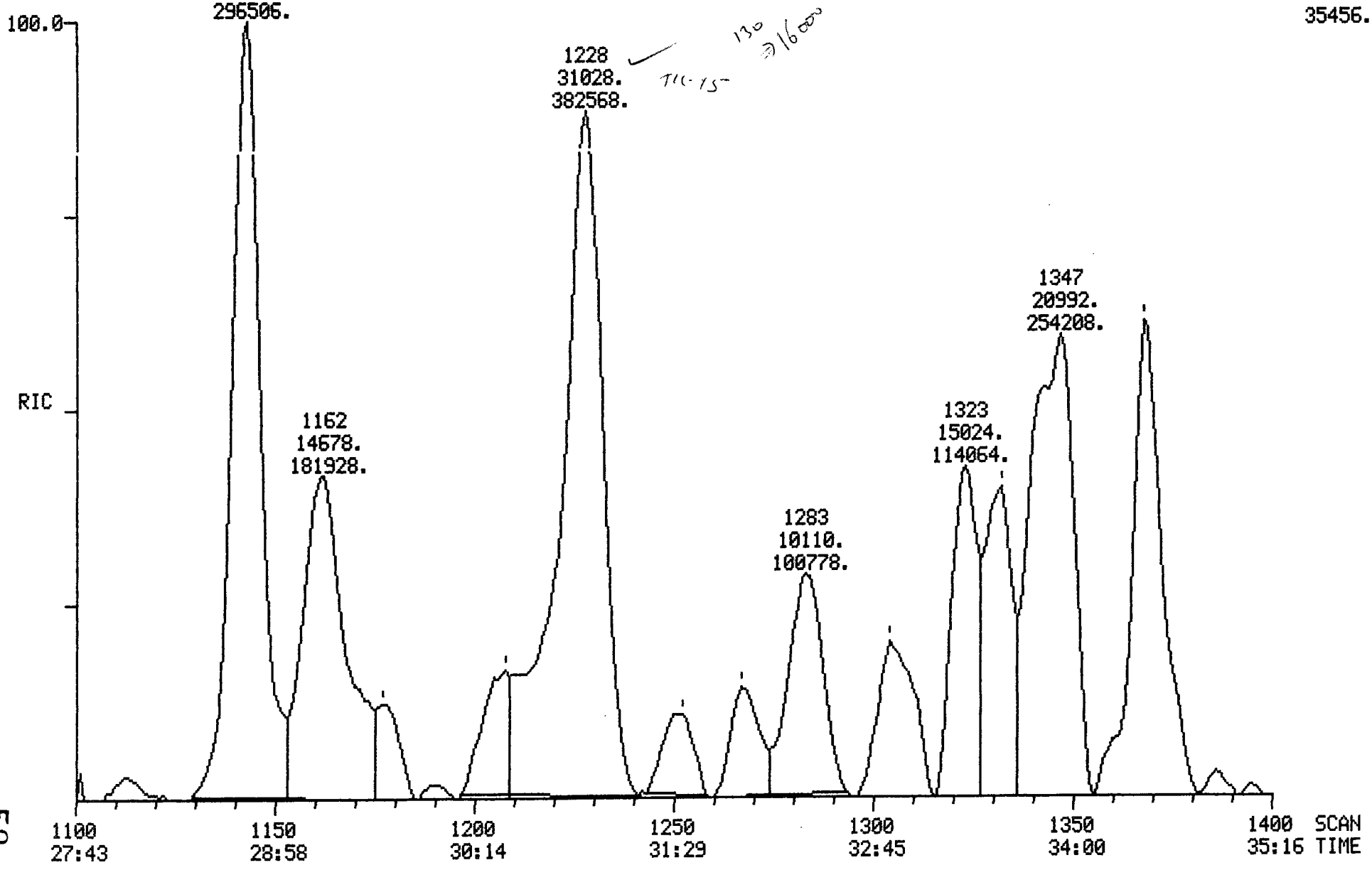
DATA: C8694 #1054
CALI: C8694 # 3

BASE M/Z: 57
RIC: 5120.



RIC
09/29/94 17:46:00
SAMPLE: 4760-001-01 1645.1 09/21/94
CONDS.: EPA METHOD 8240
RANGE: G 1,1600 LABEL: N 2, 4.0 QUAN: A 2, 1.0 J 0 BASE: U 20, 3

DATA: C8694 #1347 SCANS 1100 TO 1400
CALI: C8694 #3



Library Search Data: C8694 #1228 Base m/z: 43
09/29/94 17:46:00 + 30:56 Cali: C8694 # 3 RIC: 21920.
Sample: 4760-001-01 1645.1 09/21/94
Conds.: EPA METHOD 8240
Enhanced (S 15B 2N 0T)

42223 spectra in LIBRARYNB searched for maximum FIT
130 matched at least 8 of the 16 largest peaks in the unknown

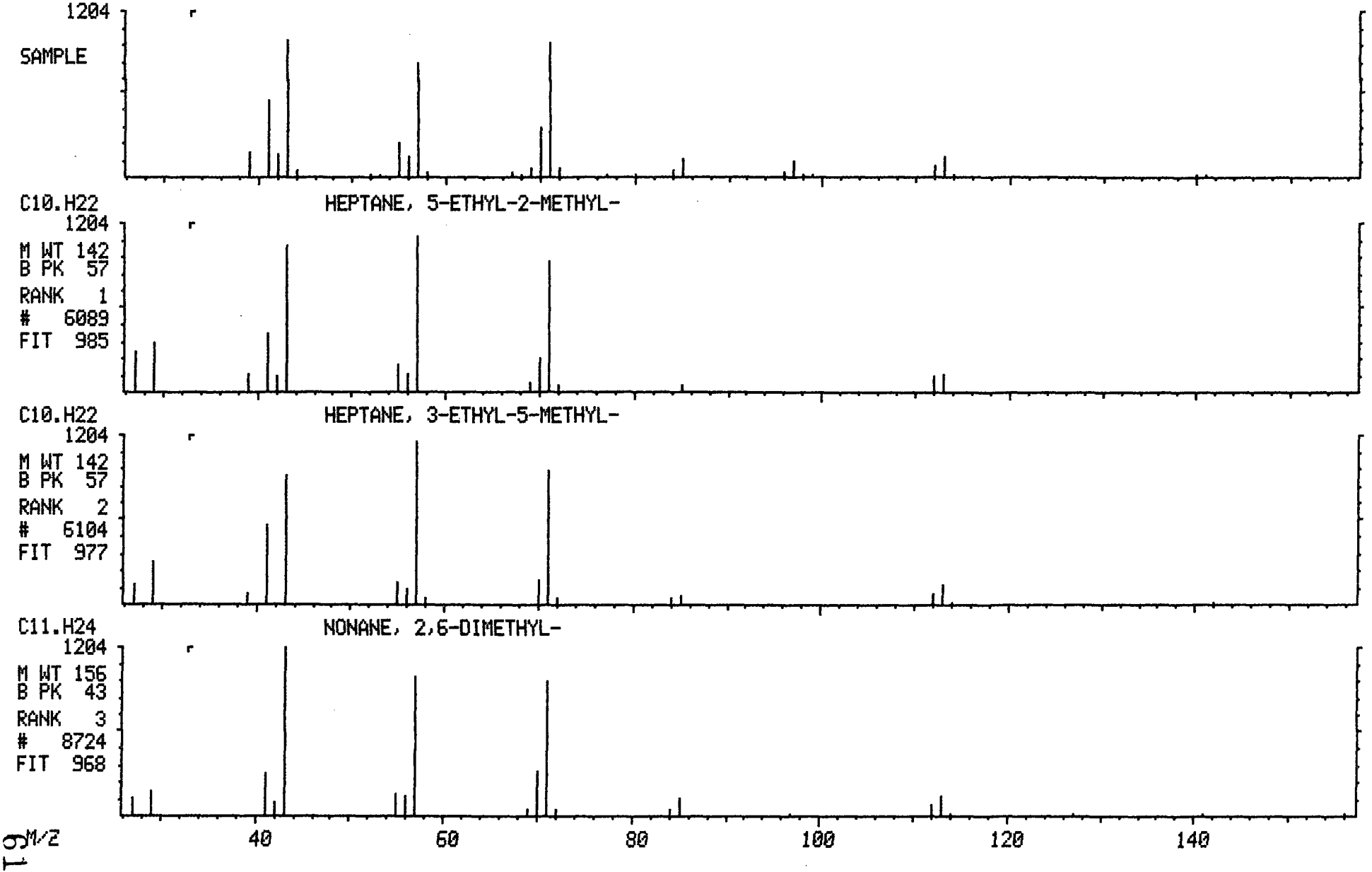
Rank	In.	Name
1	6089	HEPTANE, 5-ETHYL-2-METHYL-
2	6104	HEPTANE, 3-ETHYL-5-METHYL-
3	8724	NONANE, 2,6-DIMETHYL-
4	8732	OCTANE, 2,3,7-TRIMETHYL-
5	6085	OCTANE, 3-ETHYL-

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C10.H22	142	57	874	985	874
2	C10.H22	142	57	857	977	859
3	C11.H24	156	43	892	968	898
4	C11.H24	156	57	847	964	847
5	C10.H22	142	57	820	962	834

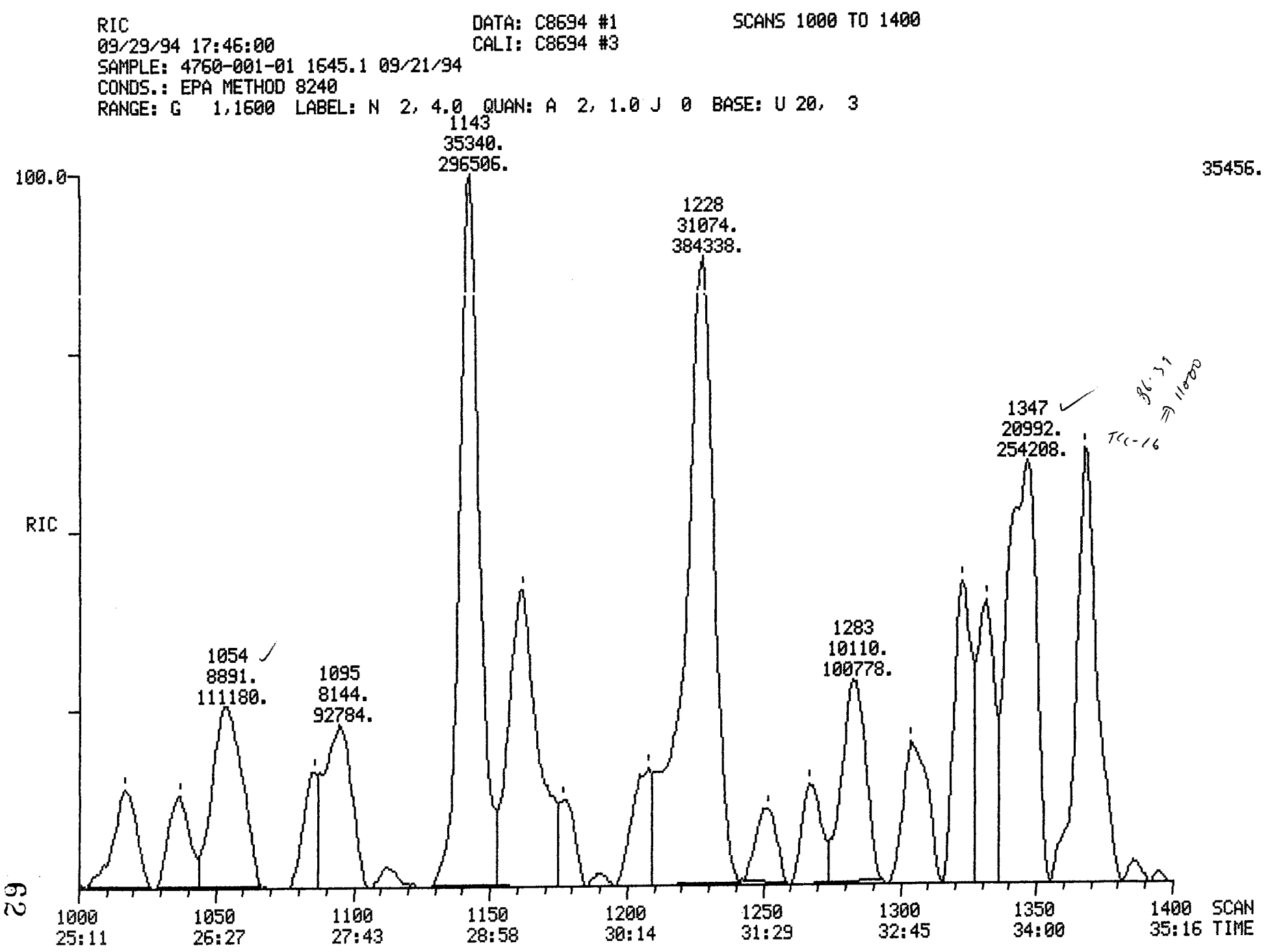
MID LIBRARY SEARCH (LIBRARYNB)
09/29/94 17:46:00 + 30:56
SAMPLE: 4760-001-01 1645.1 09/21/94
CONDS.: EPA METHOD 8240
ENHANCED (S 15B 2N 0T)

DATA: C8694 #1228
CALI: C8694 # 3

BASE M/Z: 43
RIC: 21920.



T9


RIC
09/29/94 17:46:00
SAMPLE: 4760-001-01 1645.1 09/21/94
CONDS.: EPA METHOD 8240
RANGE: G 1,1600 LABEL: N 2, 4.0 QUAN: A 2, 1.0 J 0 BASE: U 20, 3
DATA: C8694 #1
CALI: C8694 #3
SCANS 1000 TO 1400
1143
35340.
296506.
1228
31074.
384338.
1347
20992.
254208.
1054
8891.
111180.
1095
8144.
92784.
1283
10110.
100778.
35456.
100.0
RIC
62
1000
25:11
1050
26:27
1100
27:43
1150
28:58
1200
30:14
1250
31:29
1300
32:45
1350
34:00
1400
35:16
SCAN
TIME

Library Search Data: C8694 #1347 Base m/z: 43
09/29/94 17:46:00 + 33:56 Cali: C8694 # 3 RIC: 17088.
Sample: 4760-001-01 1645.1 09/21/94
Conds.: EPA METHOD 8240
Enhanced (S 15B 2N 0T)

42223 spectra in LIBRARYNB searched for maximum FIT
448 matched at least 7 of the 16 largest peaks in the unknown

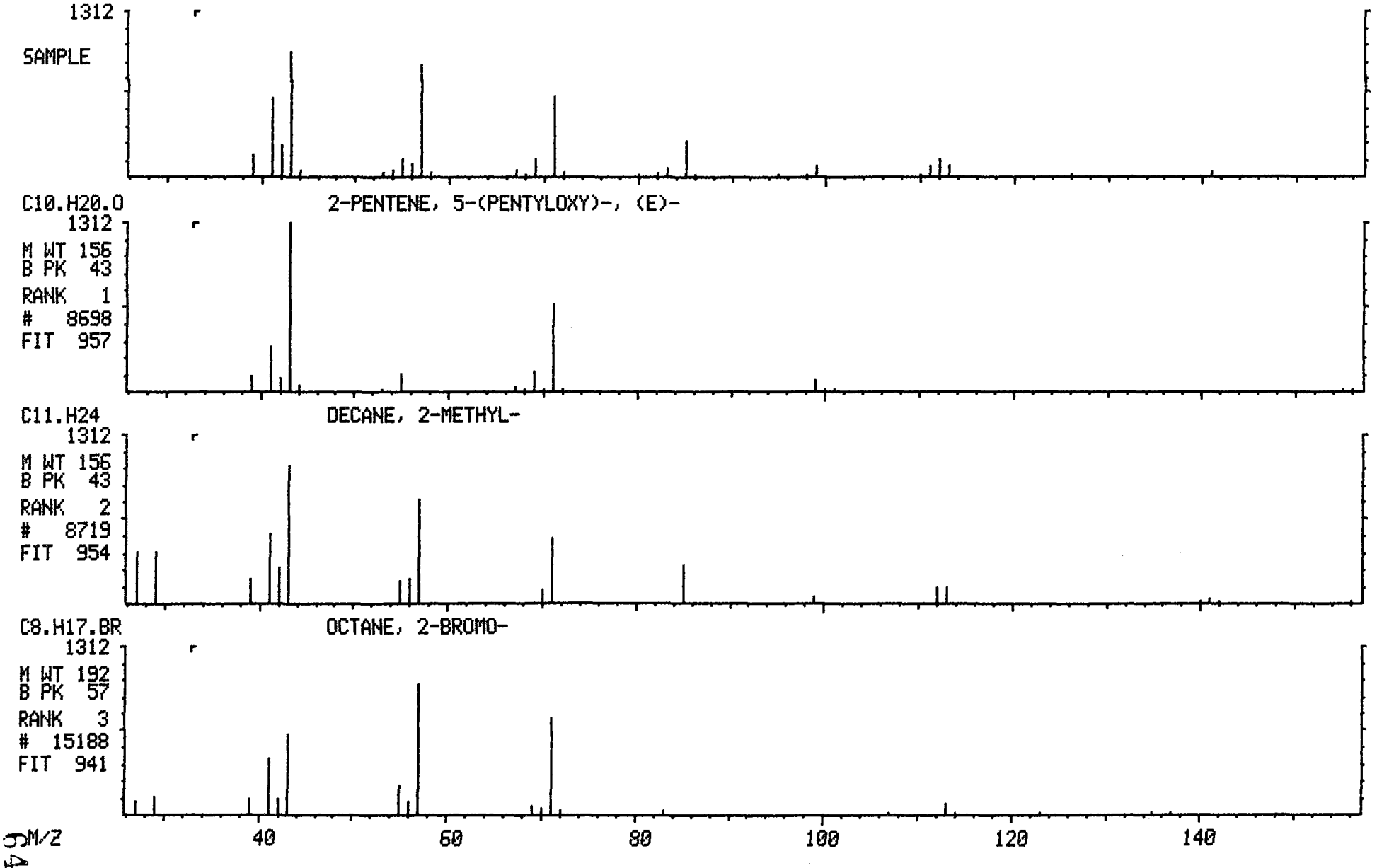
Rank	In.	Name
1	8698	2-PENTENE, 5-(PENTYLOXY)-, (E)-
2	8719	DECANE, 2-METHYL-
3	15188	OCTANE, 2-BROMO-
4	4128	HEPTANE, 2,6-DIMETHYL-
5	13877	UNDECANE, 2,10-DIMETHYL-

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C10.H20.O	156	43	516	957	524
2	C11.H24	156	43	788	954	819
3	C8.H17.BR	192	57	674	941	696
4	C9.H20	128	43	727	935	757
5	C13.H28	184	57	728	930	763

MID LIBRARY SEARCH (LIBRARYNB)
09/29/94 17:46:00 + 33:56
SAMPLE: 4760-001-01 1645.1 09/21/94
CONDS.: EPA METHOD 8240
ENHANCED (S 15B 2N 0T)

DATA: C8694 #1347
CALI: C8694 # 3

BASE M/Z: 43
RIC: 17088.



Princeton Testing Laboratory Inc.

P.O. Box 3108
3490 U.S. Route 1
Princeton, NJ 08543-3108
(609) 452-9050
(FAX) (609) 452-1959

U.S. Army, Fort Monmouth N.J.
ATTN: SELFM-PW
Building 167
Fort Monmouth, New Jersey 07703-5108
Attention: Charles Appleby

Report Date: 11/01/94
Job Number: 9404760-001
Date Received: 09/23/94
Client Job No.: 1644/1645
Page: 1

Analysis: Volatile Organics, SW, SW-846 8240
Units: ug/kg

Parameters	*Sample I.D.: 1644.1 Bldg 482 Site C-2 9/21/94 Sidewall SE*
Chloromethane	<600
Bromomethane	<600
Vinyl chloride	<600
Chloroethane	<600
Methylene chloride	470
Acetone	<300
Carbon disulfide	<300
1,1-Dichloroethene	<300
1,1-Dichloroethane	<300
1,2-Dichloroethene (Total)	<300
Chloroform	<300
1,2-Dichloroethane	<300
2-Butanone	1200
1,1,1-Trichloroethane	<300
Carbon tetrachloride	<300
Bromodichloromethane	<300
1,1,2,2-Tetrachloroethane	<300
1,2-Dichloropropane	<300
trans-1,3-Dichloropropene	<300
Trichloroethene	<300
Dibromochloromethane	<300
1,1,2-Trichloroethane	<300
Benzene	<300
cis-1,3-Dichloropropene	<300
Bromoform	<300
2-Hexanone	<300
4-Methyl-2-Pentanone	<300
Tetrachloroethene	<300
Toluene	<300
Chlorobenzene	<300
Ethylbenzene	<300
Styrene	<300
Total Xylenes	<300

RECOVERY DATA	QC LIMITS	
1,2-Dichloroethane-d4 (Surrogate)	70-121%	106
Toluene-d8 (Surrogate)	84-138%	90
4-Bromofluorobenzene (Surrogate)	59-113%	92

EPA SAMPLE NO.
1644.1 9/21

1E
VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

Lab Name: Princeton Testing Lab US_ARMY,FORT_MONMOUTH.

Lab Code: PTL **Case No.:** 4760-001 **SAS No.:** XXX **SDG No.:** XXX

Matrix: (Soil/Water) Soil

Lab Sample ID: 02

Sample wt/vol: 5 (g/mL) g

Lab File ID: C8696

Level: (low/med) med

Date Received: 09/23/94

%Moisture: not dec. 16

Date Analyzed: 09/29/94

GC Column: VOCOL ID: 0.53 mm

Dilution Factor: 50

Soil Extract Vol: ul

Soil Aliquot Vol: ul

Number TICs found: 15

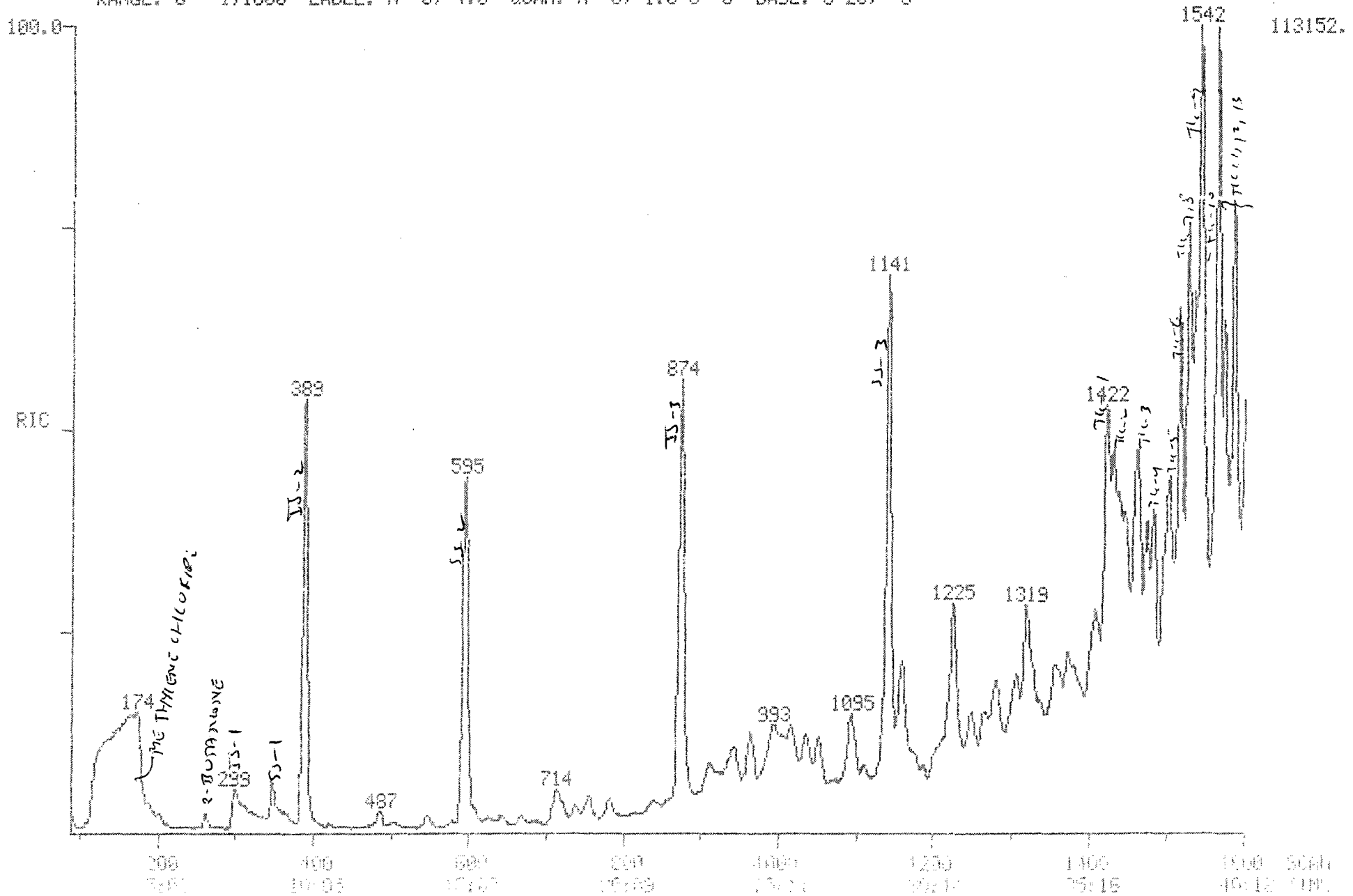
CONCENTRATION UNITS: (ug/L or ug/Kg) ug/Kg

#S	CAS NUMB	COMPOUND NAME	RT	EST. CONC.	SCAN
1	493-01-6	UNKNOWN_HYDROCARBON	35:49	1900	1422
2	29053-041	UNKNOWN_HYDROCARBON	36:01	2800	1430
3	17312-811	UNDECANE,3,5-DIMETHYL-	36:45	1400	1459
4	2980-70-3	UNKNOWN_HYDROCARBON	36:51	990	1463
5	98-06-6	BENZENE,(1,1-DIMETHYLETHYL	37:49	1100	1501
6	1758-88-9	UNKNOWN_HYDROCARBON	38:11	1300	1516
7	2958-76-1	NAPHTHALENE,DECAHYDRO-2ME	38:26	2300	1526
8	0-00-0	UNKNOWN	38:38	1800	1534
9	0-00-0	UNKNOWN	38:51	2200	1542
10	2958-76-1	NAPHTHALENE,DECAHYDRO-2ME	39:24	2800	1564
11	25155-151	BENZENE,METHYL(1-METHYLETH	39:37	1100	1573
12	527-53-7	BENZENE,1,2,3,5-TETRAMETHYL	39:56	1700	1585
13	0-00-0	UNKNOWN	39:59	-750	1587
14	0-00-0	UNKNOWN_HYDROCARBON	30:51	2100	1225
15	0-00-0	UNKNOWN_HYDROCARBON	33:14	1900	1319

RIC
09/29/94 19:28:00
SAMPLE: 4760-001-02 1644.1 09/21/94
CONDS.: EPA METHOD 8240
RANGE: G 1,1600 LABEL: N 0, 4.0 QUAN: A 0, 1.0 J 0 BASE: U 20, 3

DATA: C8696 #1
CALI: C8696 #3

SCANS 88 TO 1600



Quantitation Report File: C8696

Data: C8696.TI
9/29/94 19:28:00
Sample: 4760-001-02 1644.1 09/21/94
Conds.: EPA METHOD 8240
Formula: X50DIL Instrument: FINN Weight: 0.000
Submitted by: USARMY Analyst: UC Acct. No.: 4760-001

AMOUNT=AREA * REF AMNT/(REF AREA * RESP FACT)
Resp. fac. from Library Entry

No	Name
1	C101 BROMOCHLOROMETHANE **INT. STD. **
2	C110 1,4-DIFLUOROBENZENE **INT. STD. **
3	C120 CHLOROBENZENE-D5 **INT. STD. **
4	CS15 1,2-DICHLOROETHANE-D4 **S. STD. **
5	CS05 TOLUENE-D8 **S. STD. **
6	CS10 4-BROMOFLUOROBENZENE **S. STD. **
7	C010 CHLOROMETHANE **
8	C015 BROMOMETHANE
9	C020 VINYL CHLORIDE *
10	C025 CHLOROETHANE
11	C030 METHYLENE CHLORIDE
12	C251 ACROLIN
13	C035 ACETONE
14	C252 ACRYLONITRILE
15	C040 CARBON DISULFIDE
16	C045 1,1-DICHLOROETHENE *
17	C050 1,1-DICHLOROETHANE **
18	C055 TRANS-1,2-DICHLOROETHENE
19	C000 TRICHLOROFLUOROMETHANE
20	C060 CHLOROFORM *
21	C065 1,2-DICHLOROETHANE
22	C110 2-BUTANONE
23	C115 1,1,1-TRICHLOROETHANE
24	C120 CARBON TETRACHLORIDE
25	C125 VINYL ACETATE
26	C130 BROMO DICHLOROMETHANE
27	C140 1,2-DICHLOROPROPANE *
28	C145 TRANS-1,3 DICHLOROPROPENE
29	C150 TRICHLOROETHENE
30	C155 DIBROMOCHLOROMETHANE
31	C160 1,1,2-TRICHLOROETHANE
32	C165 BENZENE
33	C143 CIS-1,3-DICHLOROPROPENE
34	C175 2-CHLOROETHYL VINYL ETHER
35	C180 BROMOFORM **
36	C220 TETRACHLOROETHENE
37	C210 2-HEXANONE
38	C205 4-METHYL 2-PENTANONE
39	C225 1,1,2,2-TETRACHLOROETHANE **
40	C230 TOLUENE *
41	C235 CHLOROBENZENE **
42	C240 ETHYL BENZENE *
43	C245 STYRENE
44	C250 M+P-XYLENES
45	C253 1,3-DICHLOROBENZENE
46	C254 1,4-DICHLOROBENZENE
47	C255 1,2-DICHLOROBENZENE

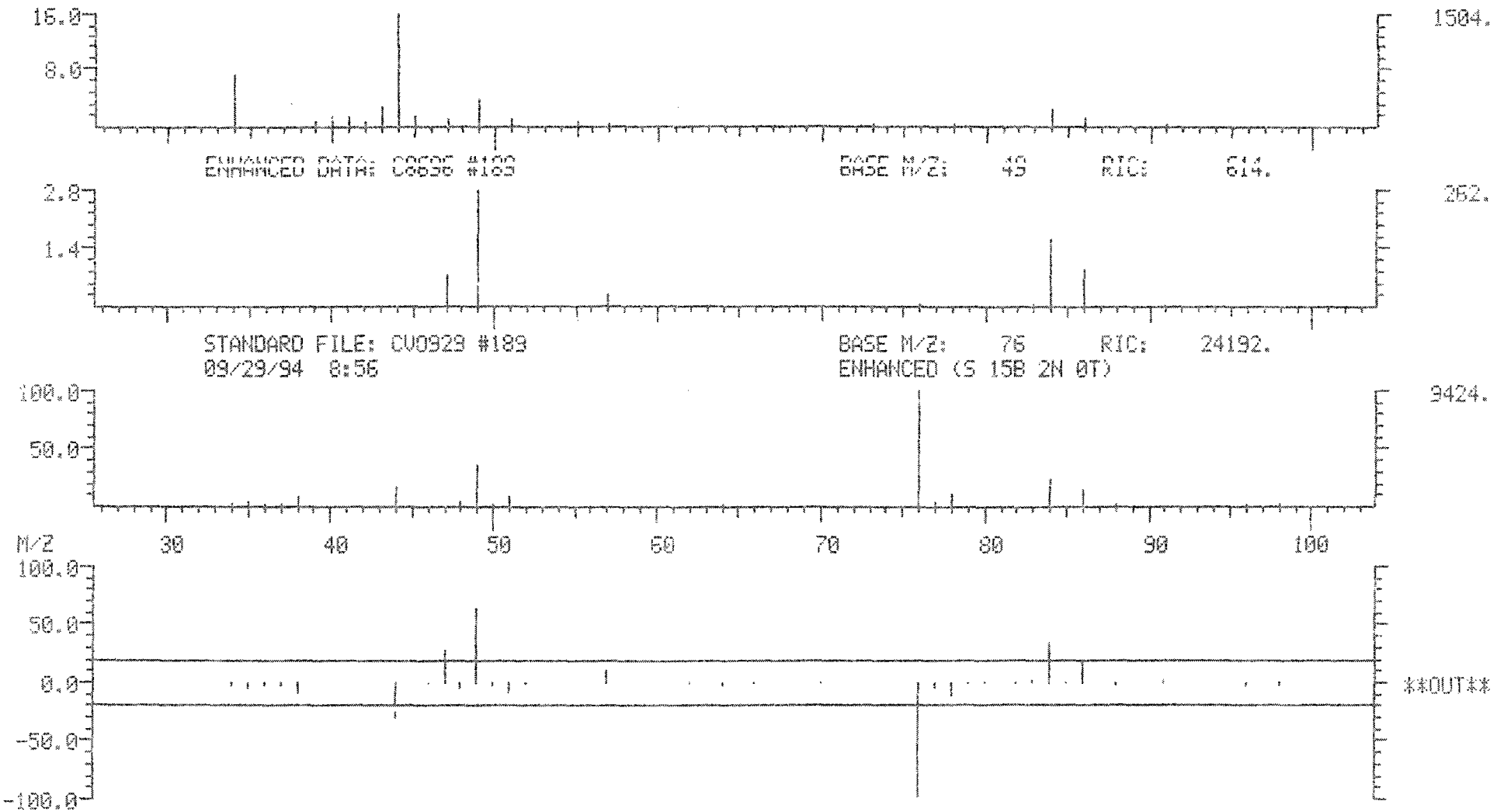
CB696

No	Name
46	C250 O-XYLENE
45	BENZENE-D6 **S. STD. **

No	m/z	Scan	Time	Ref	RRT	Meth	Area(Hght)	Amount	%Tot	
1	49	299	7:32	1	1.000	A BB	8235.	50.000 NG	15.51	
2	114	389	9:48	2	1.000	A BB	144726.	50.000 NG	15.51	
3	117	873	21:59	3	1.000	A BB	114280.	50.000 NG	15.51	
4	65	347	8:44	1	1.161	M XX	14700.	52.773 NG	16.37	106
5	98	595	14:59	3	0.682	A BB	138340.	45.169 NG	14.01	90
6	95	1141	28:45	3	1.307	A BB	101483.	45.857 NG	14.23	92
7	NOT FOUND									
8	NOT FOUND									
9	NOT FOUND									
10	NOT FOUND									
11	49	189	4:46	1	0.632	A BB	3159.	7.854 NG	2.44	
12	NOT FOUND									
13	NOT FOUND									
14	NOT FOUND									
15	NOT FOUND									
16	NOT FOUND									
17	NOT FOUND									
18	NOT FOUND									
19	NOT FOUND									
20	NOT FOUND									
21	NOT FOUND									
22	43	261	6:34	2	0.671	A BB	8965.	20.258 NG	6.28	
23	NOT FOUND									
24	NOT FOUND									
25	NOT FOUND									
26	NOT FOUND									
27	NOT FOUND									
28	NOT FOUND									
29	NOT FOUND									
30	NOT FOUND									
31	NOT FOUND									
32	NOT FOUND									
33	NOT FOUND									
34	NOT FOUND									
35	NOT FOUND									
36	NOT FOUND									
37	NOT FOUND									
38	NOT FOUND									
39	NOT FOUND									
40	91	607	15:17	3	0.695	A BB	1971.	0.~~557~~ NG	0.17	
41	NOT FOUND									
42	106	914	23:01	3	1.047	A BB	34.	0.~~035~~ NG	0.01	
43	NOT FOUND									
44	NOT FOUND									
45	NOT FOUND									
46	NOT FOUND									
47	NOT FOUND									
48	NOT FOUND									
*	NOT FOUND									

DATA FILE: C8696 #189 STANDARD FILE: CV0929 #189
TARGET COMPOUND COMPARISON CALI: C8696 #3
COMPOUND: C030 METHYLENE CHLORIDE

RAW DATA: C8696 #189 BASE M/Z: 44 RIC: 4044.
09/29/94 19:28

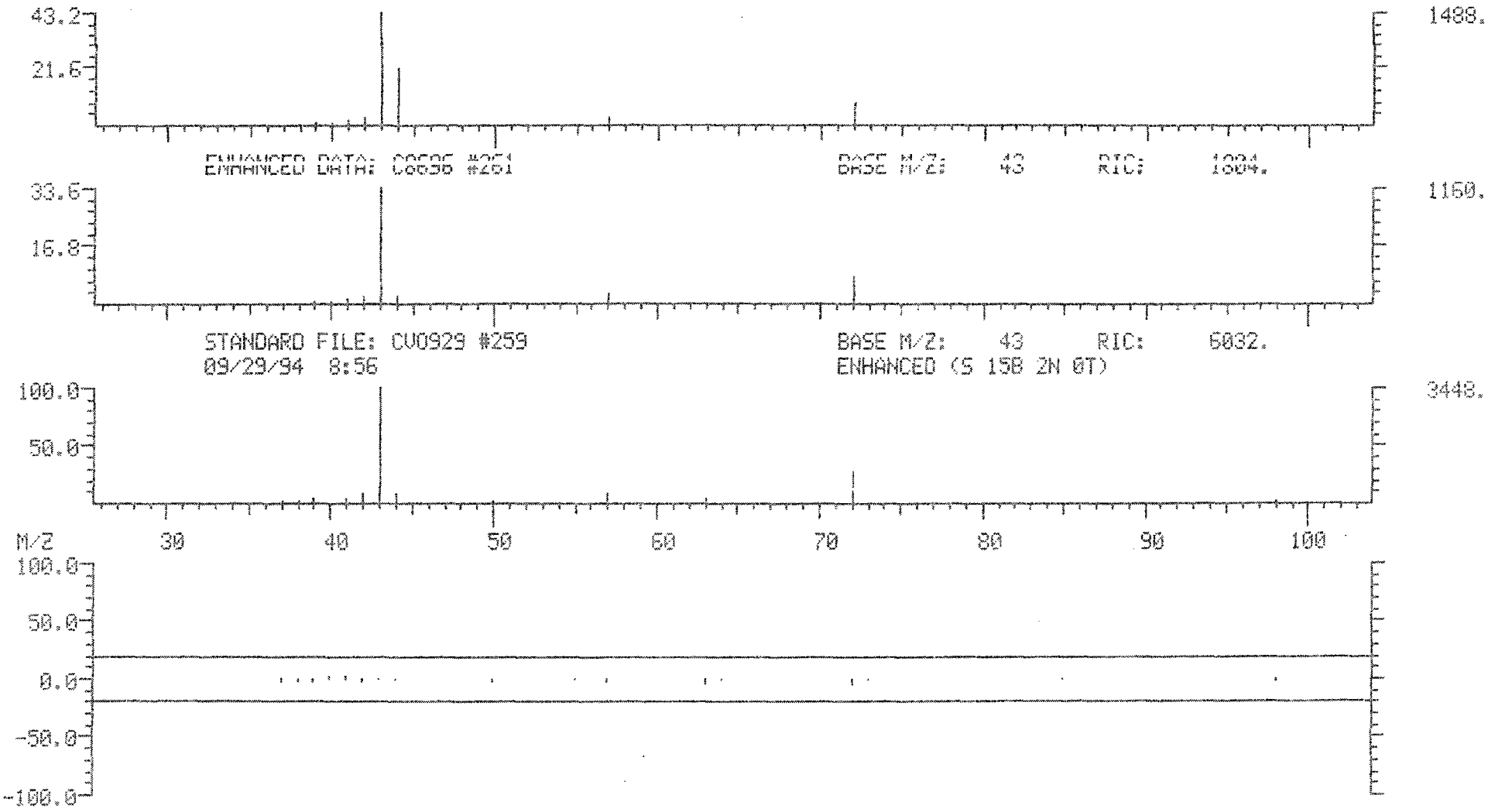


DATA FILE: C8696 #261
TARGET COMPOUND COMPARISON
COMPOUND: C110 2-BUTANONE

STANDARD FILE: CV0929 #259
CALI: C8696 #3

RAW DATA: C8696 #261
09/29/94 19:28

BASE M/Z: 43 RIC: 2796.



PROCEDURE: FILTER/TIC DIAGNOSTIC REPORT 9/29/94 20:36:22

DATA FILE: C8696

FILTER SCAN PARAMETERS

MAX NUMBER TICS:	15
MAX TABLE ENTRIES:	539
SCAN TOLERANCE :	2
MIN RIC HT. [%]:	10
FIRST SCAN :	1
LAST SCAN :	1600
TIC THRESHOLD :	600

METHOD LIBRARY & LISTS

TIC I.S. LIBRARY:	LIBRARYLS
NBS SEARCH PROC :	SERLIB
PEAK FINDER PROC:	VOME
TCA I.S. LL :	LS
FILE NAME LIST :	TCAREF2

TARGET COMPOUND ANALYSIS:

TARGETS (QUAN LIST)	IS PEAKS	TOTAL TARGET PEAKS
7	3	10

FILTER PROCESSING:

	<-----REJECT PEAKS----->						
TOTAL PEAKS	< 1ST SCAN	> LAST SCAN	< MIN RIC HT	< SCAN TOL	> MAX # PEAKS	TOTAL REJECTS	TOTAL TICS
31	0	0	1	6	9	16	15

TIC PROCESSING:

ID	SCAN#	PURITY	FIT	MW	COMPOUND NAME [BEFORE TIC THRESHOLD]
1	83	408	426	85	ACETIC ACID, CYANO-
2	1422	834	965	138	NAPHTHALENE, DECAHYDRO-, CIS-
3	1431	639	861	140	CYCLOPENTANE, 1-METHYL-3-(2-ME
4	1459	761	906	184	UNDECANE, 3,5-DIMETHYL-
5	1463	683	854	194	CYCLOPENTANE, 1,1'-(1,4-BUTAND
6	1501	777	952	134	BENZENE, (1,1-DIMETHYLETHYL)-
7	1516	753	969	134	BENZENE, 2-ETHYL-1,4-DIMETHYL-
8	1526	786	989	152	NAPHTHALENE, DECAHYDRO-2-METHY
9	1534	487	698	247	AZIRIDINONE, 1-(1,1-DIMETHYLET
10	1542	449	878	238	CYCLOHEXANE, UNDECYL-
11	1562	624	992	152	NAPHTHALENE, DECAHYDRO-2-METHY
12	1564	651	986	152	NAPHTHALENE, DECAHYDRO-2-METHY
13	1573	715	917	134	BENZENE, METHYL(1-METHYLETHYL)
14	1585	741	979	134	BENZENE, 1,2,3,5-TETRAMETHYL-
15	1587	500	909	184	UNDECANE, 4,8-DIMETHYL-

Quantitation Report File: C8696

Data: C8696.TI
09/29/94 19:28:00
Sample: 4760-001-02 1644.1 09/21/94
Comas.: EPA METHOD 8240
Formula. X50DIL Instrument: FINN Weight: 0.000
Submitted by: USARMY Analyst: UC Acct. No.: 4760-001

AMOUNT=AREA * REF AMNT/(REF AREA * RESP FACT)
Resp. fac. from Library Entry

No	CAS #	Name
1	0-00-0	CI01 BROMOCHLOROMETHANE **INT. STD. **
2	0-00-0	CI10 1,4-DIFLUOROBENZENE **INT. STD. **
3	0-00-0	CI20 CHLOROBENZENE-D5 **INT. STD. **
4	0-00-0	UNKNOWN
5	493-01-6	NAPHTHALENE, DECAHYDRO-, CIS-
6	29053-04-1	CYCLOPENTANE, 1-METHYL-3-(2-METHYLPROPYL)-
7	17312-81-1	UNDECANE, 3,5-DIMETHYL-
8	2980-70-3	CYCLOPENTANE, 1,1'-(1,4-BUTANDIYL)BIS-
9	98-06-6	BENZENE, (1,1-DIMETHYLETHYL)-
10	1758-88-9	BENZENE, 2-ETHYL-1,4-DIMETHYL-
11	2958-76-1	NAPHTHALENE, DECAHYDRO-2-METHYL-
12	0-00-0	UNKNOWN
13	0-00-0	UNKNOWN
14	2958-76-1	NAPHTHALENE, DECAHYDRO-2-METHYL-
~~15~~	~~2958-76-1~~	~~NAPHTHALENE, DECAHYDRO-2-METHYL-~~
16	25155-15-1	BENZENE, METHYL(1-METHYLETHYL)-
17	527-53-7	BENZENE, 1,2,3,5-TETRAMETHYL-
18	0-00-0	UNKNOWN

No	m/z	Scan	Time	Ref	RRT	Meth	Area(Hght)	Amount	%Tot
1	TOT	299	7:32	0	ISINV	A BB	51988.	****** UG/L	00.00
2	TOT	389	9:48	0	ISINV	A BB	319250.	****** UG/L	00.00
3	TOT	873	21:59	0	ISINV	A BB	299992.	****** UG/L	00.00
4	TOT	83	2:05	1	0.278	A BB	621550.	~~597.787~~	58.89
5	TOT	1422	35:49	3	1.629	A BV	189679.	31.614 1900	3.11
6	TOT	1430	36:01	3	1.638	A VB	282162.	47.028 2800	4.63
7	TOT	1459	36:45	3	1.671	A BV	137024.	22.838 1400	2.25
8	TOT	1463	36:51	3	1.676	A BV	99619.	16.604 990	1.64
9	TOT	1501	37:49	3	1.719	A BB	109252.	18.209 1100	1.79
10	TOT	1516	38:11	3	1.737	A BB	125960.	20.994 1300	2.07
11	TOT	1526	38:26	3	1.748	A BV	230368.	38.396 2300	3.78
12	TOT	1534	38:38	3	1.757	A BV	178871.	29.813 1800	2.94
13	TOT	1542	38:51	3	1.766	A VB	219472.	36.580 2200	3.60
14	TOT	1564	39:24	3	1.792	A BV	284616.	47.437 2800	4.67
~~15~~	~~TOT~~	~~1564~~	~~39:24~~	~~3~~	~~1.792~~	~~A BV~~	~~284616.~~	47.437	4.67
16	TOT	1573	39:37	3	1.802	A VV	115508.	19.252 1100	1.90
17	TOT	1585	39:56	3	1.816	A VB	171501.	28.584 1700	2.82
18	TOT	1587	39:59	3	1.818	A BB	75168.	12.528 1250	1.23

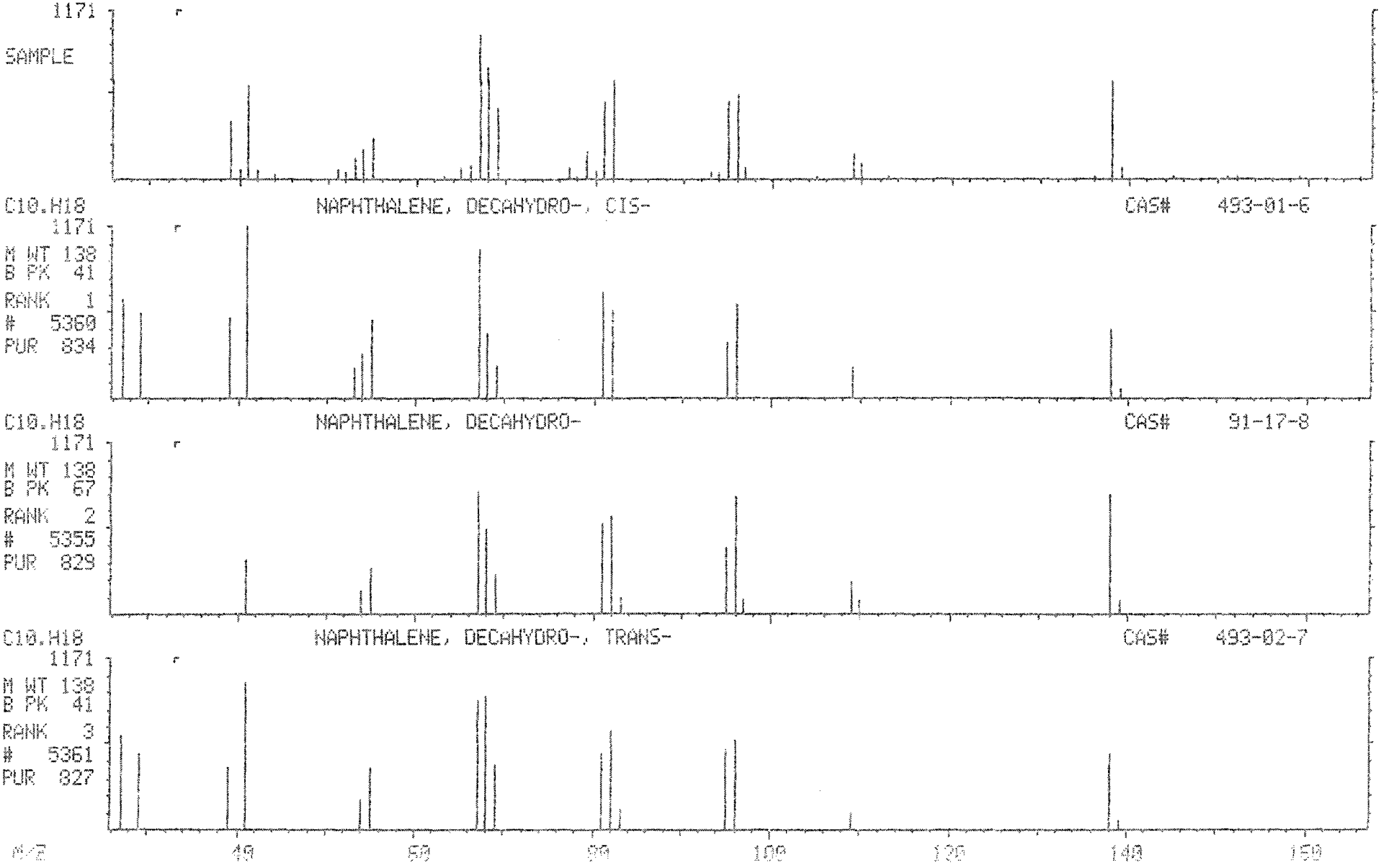
No	Ret(L)	Ratio	RRT(L)	Ratio	Amnt	Amnt(L)	R.Fac	R.Fac(L)	Ratio
1	9:13	0.82	1.000						
2	19:21	0.51	1.000						
3	23:54	0.92	1.000						

No	Ret(L) Ratio	RRT(L) Ratio	Amnt	Amnt(L)	R.Fac	R.Fac(L)	Ratio
4			597.79	1.00	597.787	1.000	597.79
5			31.61	1.00	31.614	1.000	31.61
6			47.03	1.00	47.028	1.000	47.03
7			22.84	1.00	22.838	1.000	22.84
8			16.60	1.00	16.604	1.000	16.60
9			18.21	1.00	18.209	1.000	18.21
10			20.99	1.00	20.994	1.000	20.99
11			38.40	1.00	38.396	1.000	38.40
12			29.81	1.00	29.813	1.000	29.81
13			36.58	1.00	36.580	1.000	36.58
14			47.44	1.00	47.437	1.000	47.44
15			47.44	1.00	47.437	1.000	47.44
16			19.25	1.00	19.252	1.000	19.25
17			28.58	1.00	28.584	1.000	28.58
18			12.53	1.00	12.528	1.000	12.53

MID LIBRARY SEARCH (LIBRARYNB)
09/29/94 19:28:00 + 35:49
SAMPLE: 4760-001-02 1644.1 09/21/94
CONDS.: EPA METHOD 8240
ENHANCED (S 15B 2N 0T)

DATA: C8696 #1422
CALI: C8696 # 3

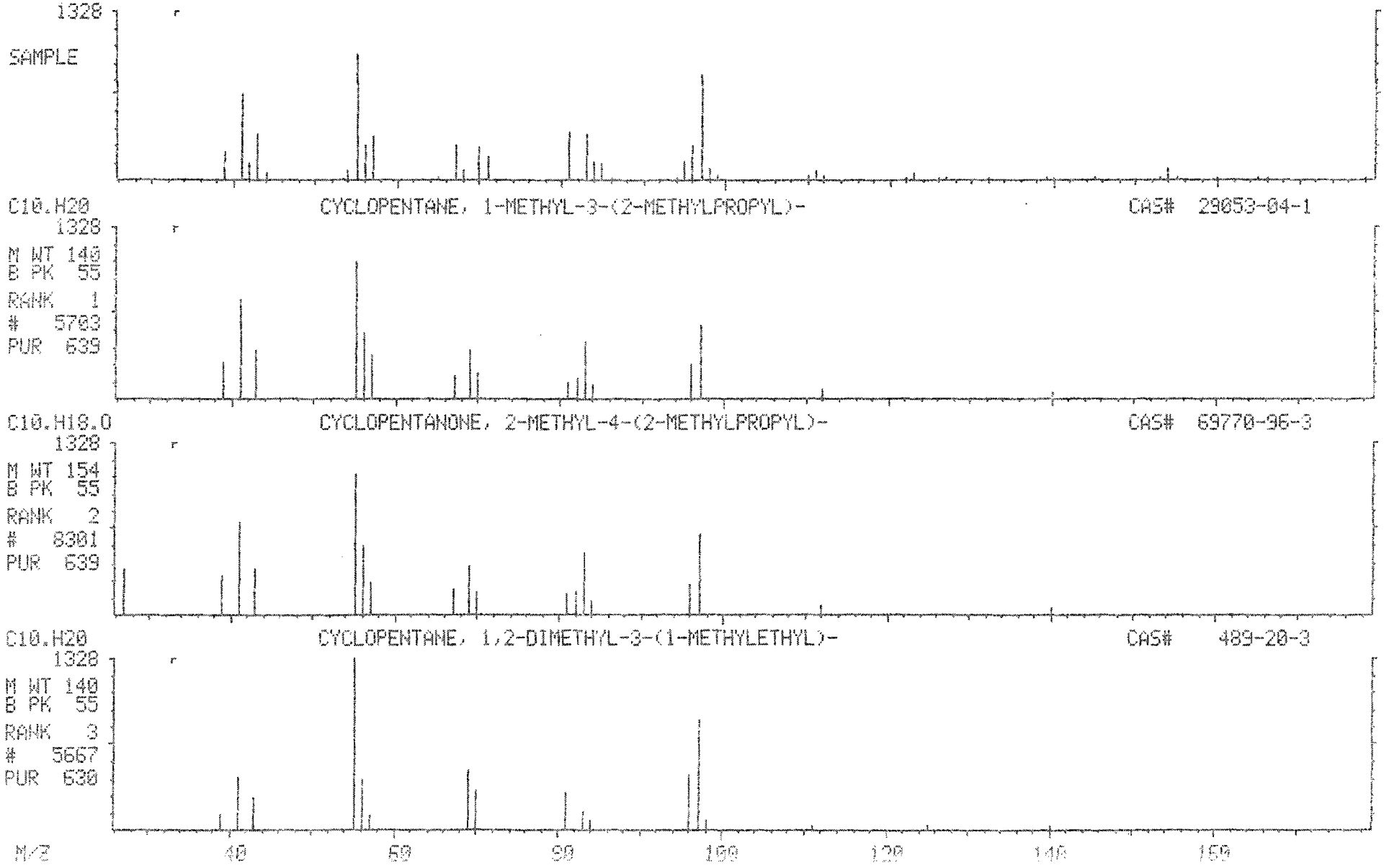
BASE M/Z: 67
RIC: 22016.



MID LIBRARY SEARCH (LIBRARYNB)
09/29/94 19:28:00 + 35:03
SAMPLE: 4760-001-02 1644.1 09/21/94
CONDS.: EPA METHOD 8240
ENHANCED (S 15B 2N 0T)

DATA: C8696 #1431
CALI: C8696 # 3

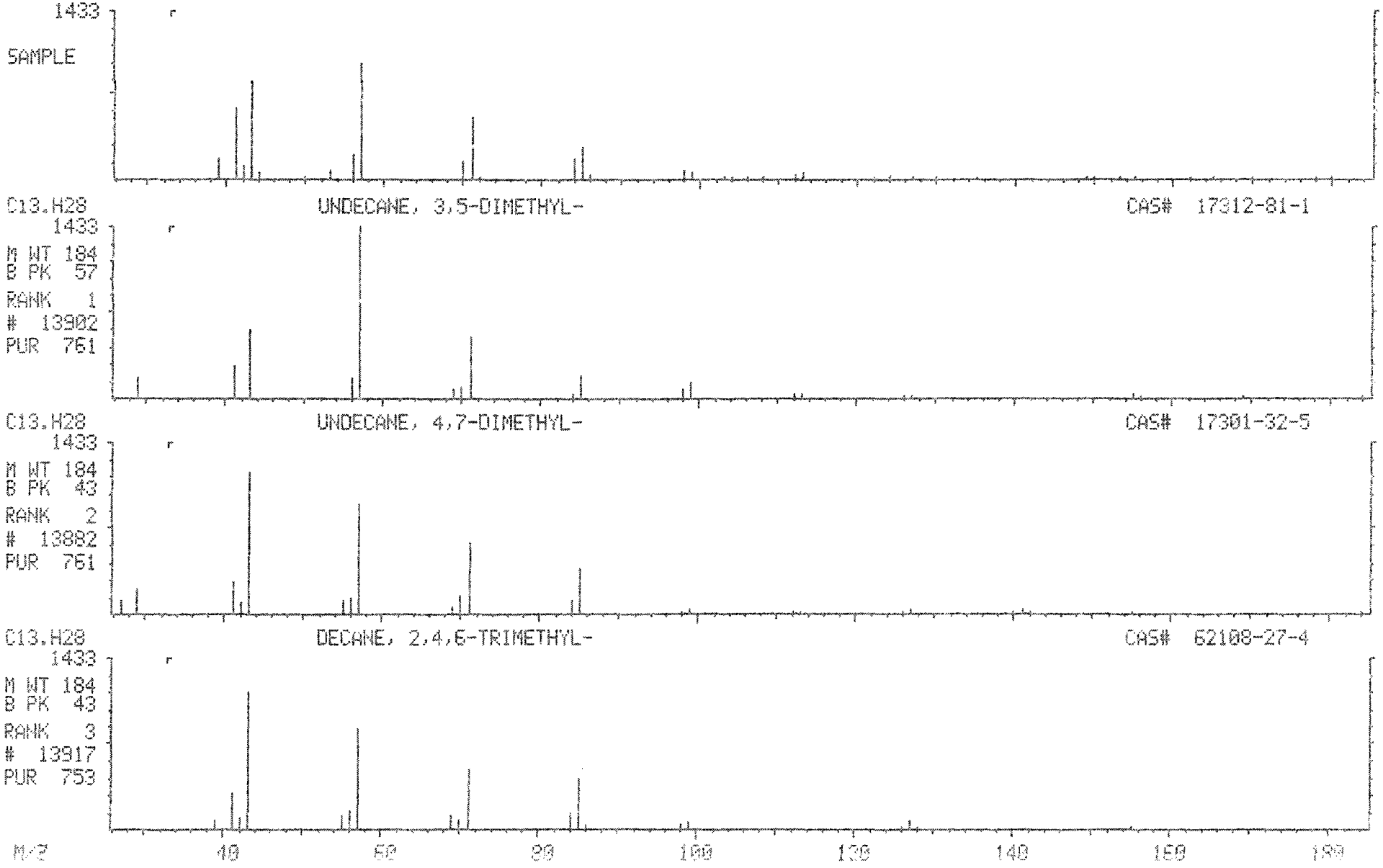
BASE M/Z: 55
RIC: 17088.



MID LIBRARY SEARCH (LIBRARYNB)
09/29/94 19:28:00 + 36:45
SAMPLE: 4760-001-02 1644.1 09/21/94
CONDS.: EPA METHOD 8240
ENHANCED (S 15B 2N 0T)

DATA: C8696 #1459
CALI: C8696 # 3

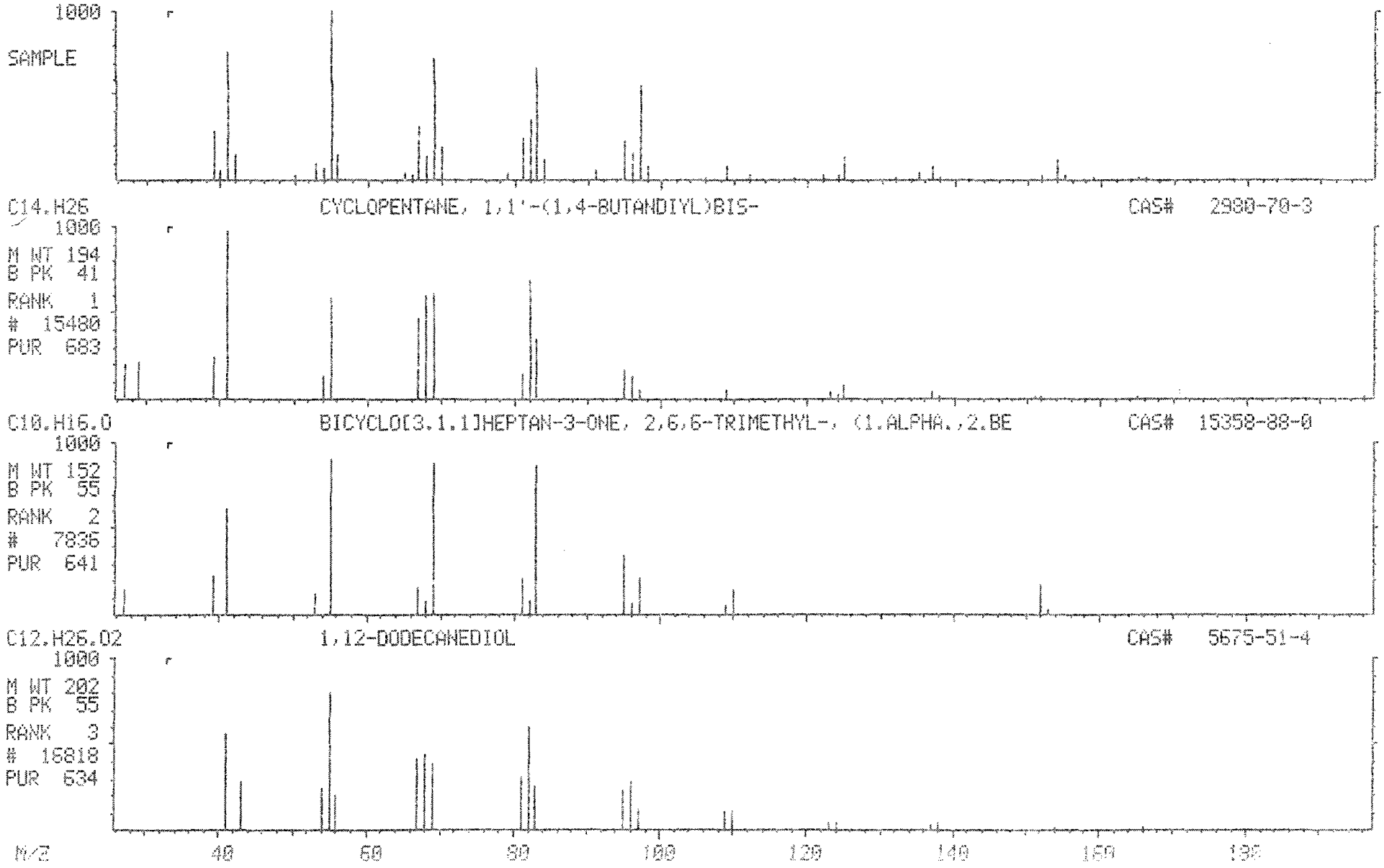
BASE M/Z: 57
RIC: 16608.



MID LIBRARY SEARCH (LIBRARYNB)
09/29/94 19:28:00 + 36:51
SAMPLE: 4760-001-02 1644.1 09/21/94
CONDS.: EPA METHOD 8240
ENHANCED (S 15B 2N 0T)

DATA: C8696 #1463
CALI: C8696 # 3

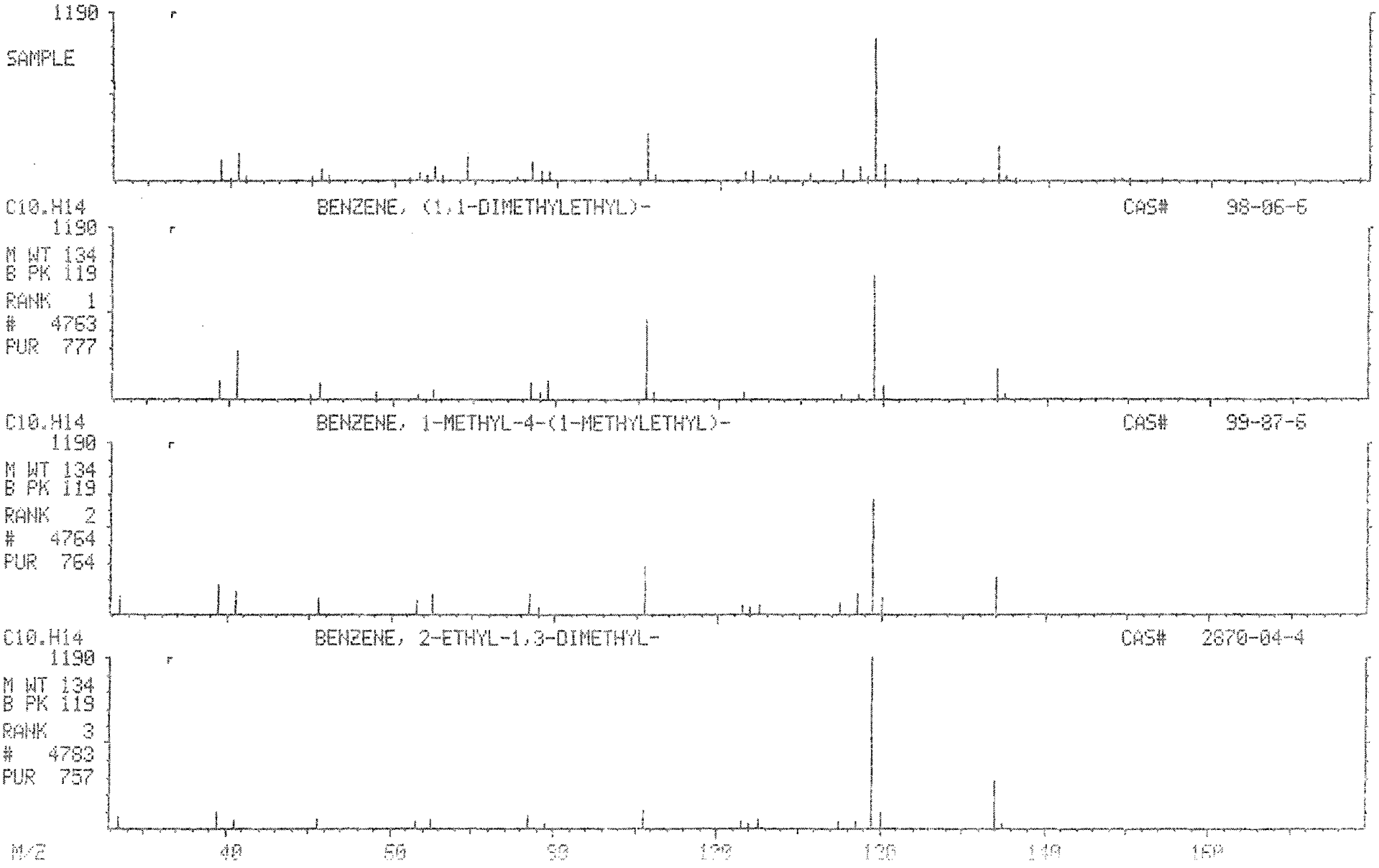
BASE M/Z: 55
RIC: 16992.



MID LIBRARY SEARCH (LIBRARYNB)
09/29/94 19:28:00 + 37:49
SAMPLE: 4760-001-02 1644.1 09/21/94
CONDS.: EPA METHOD 8240
ENHANCED (S 15B 2N 0T)

DATA: C8696 #1501
CALI: C8696 # 3

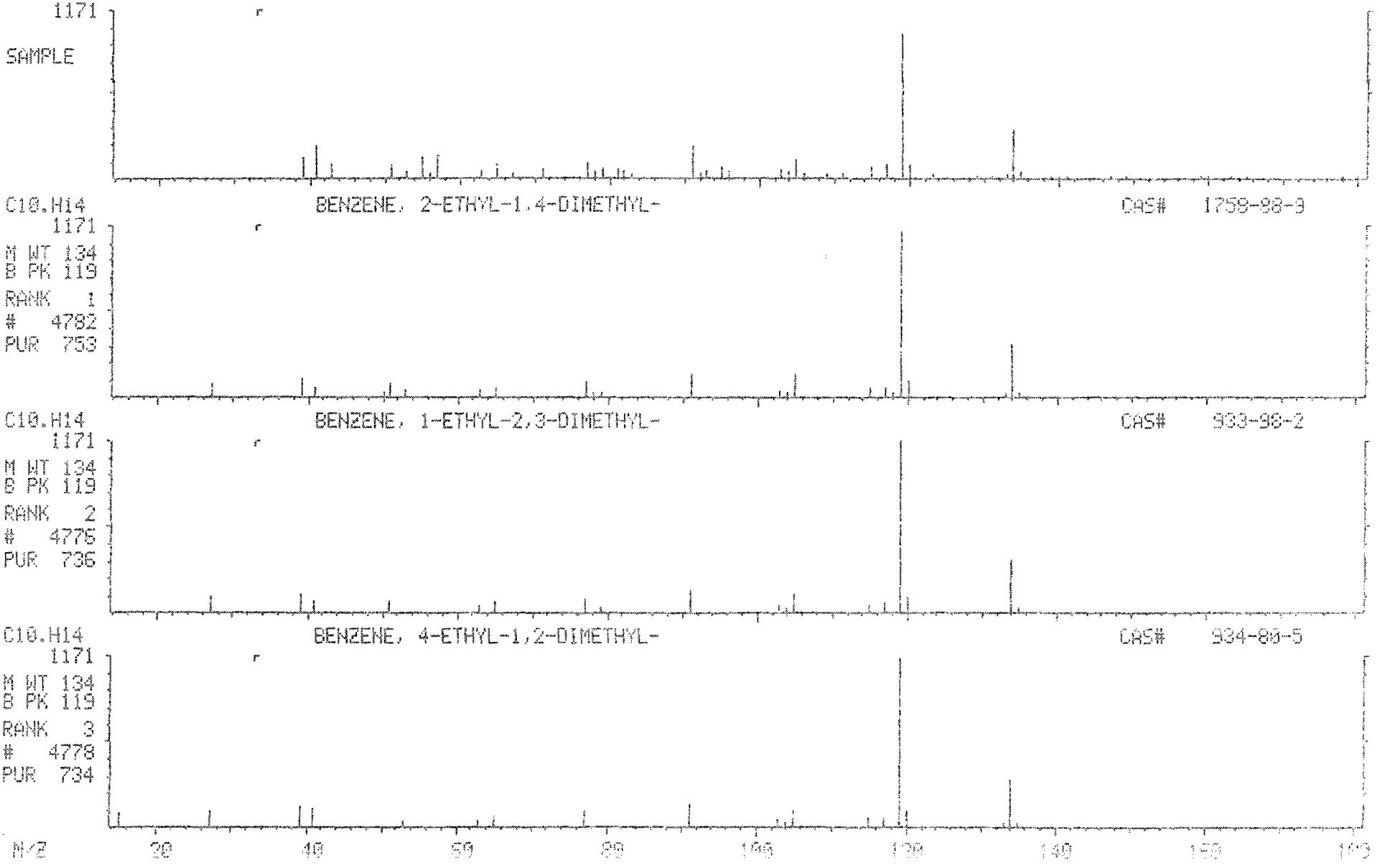
BASE M/Z: 119
RIC: 13920.



MID LIBRARY SEARCH (LIBRARYNB)
09/29/94 19:28:00 + 38:11
SAMPLE: 4760-001-02 1644.1 09/21/94
CONDS.: EPA METHOD 8240
ENHANCED (S 15B 2N 0T)

DATA: C8696 #1516
CALI: C8696 # 3

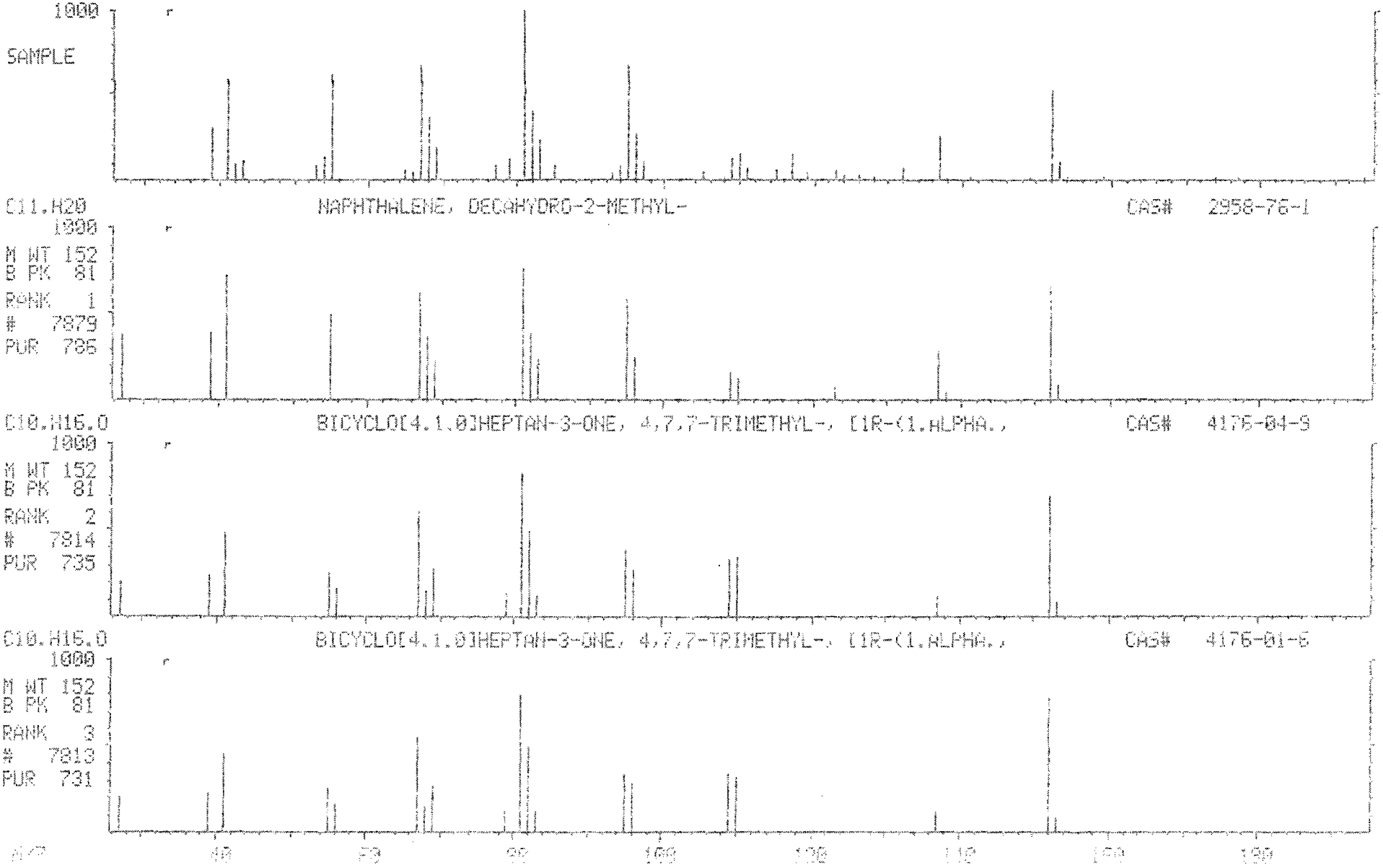
BASE M/Z: 119
RIC: 30816.



MID LIBRARY SEARCH (LIBRARYNB)
09/29/94 19:28:00 + 38:26
SAMPLE: 4760-001-02 1644.1 09/21/94
CONDS.: EPA METHOD 8240
ENHANCED (S 15B 2N 0T)

DATA: C8696 #1525
CALI: C8696 # 3

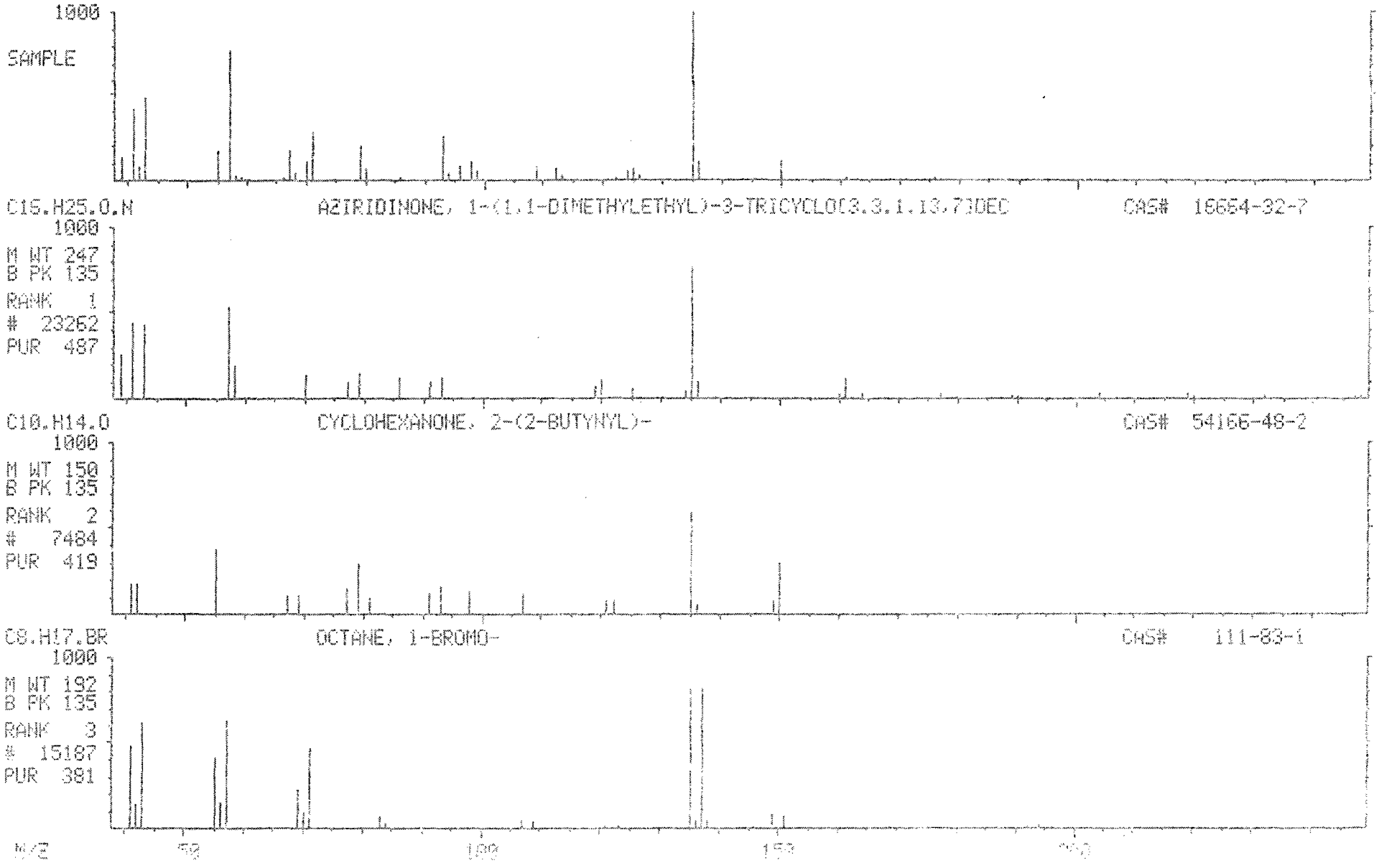
BASE M/Z: 81
RIC: 34304.



MID LIBRARY SEARCH (LIBRARYN8)
09/29/94 19:28:00 + 38:38
SAMPLE: 4760-001-02 1644.1 09/21/94
CONDS.: EPA METHOD 8240
ENHANCED (S 15B 2N 0T)

DATA: C8696 #1534
CALI: C8696 # 3

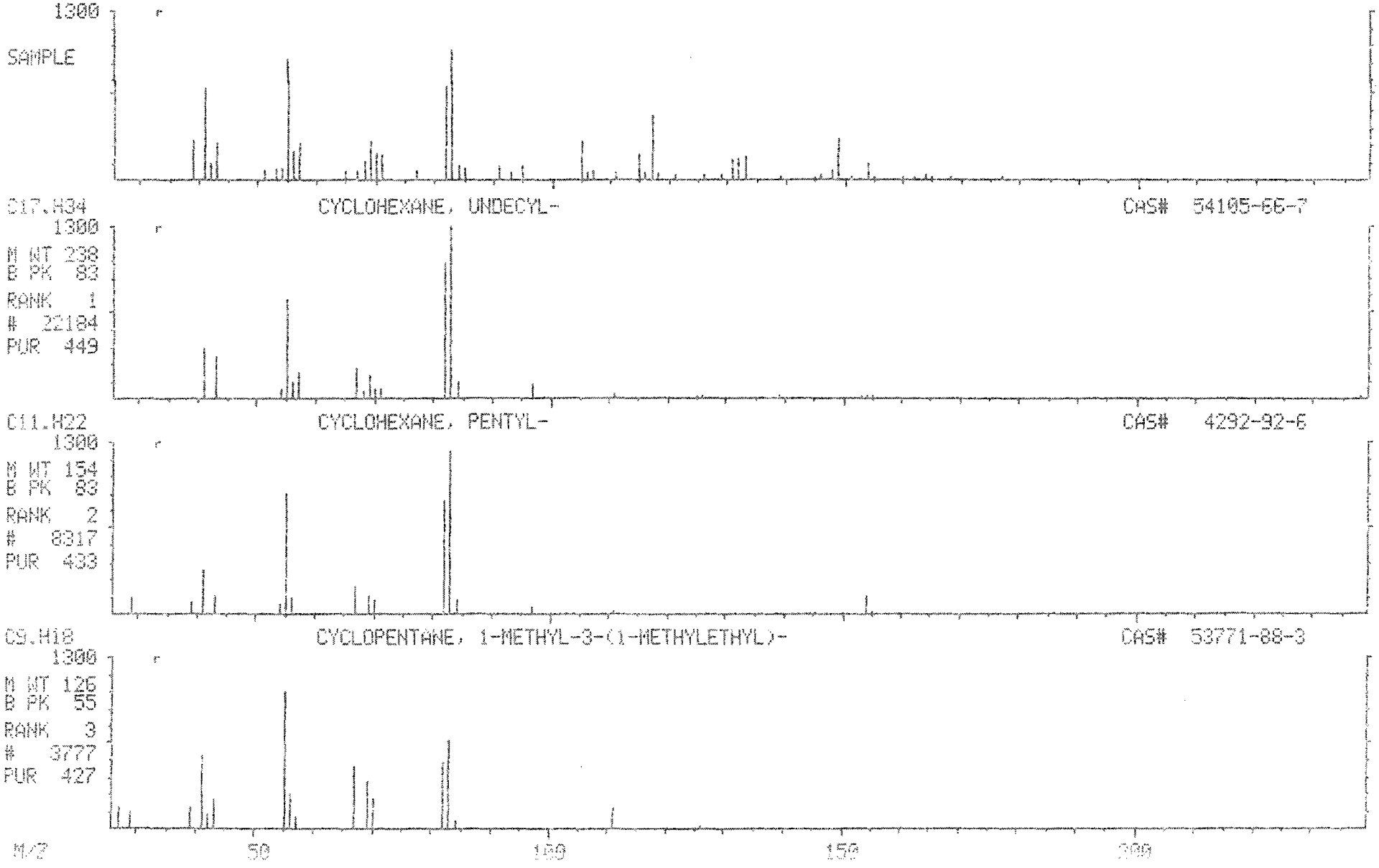
BASE M/Z: 135
RIC: 17664.



MID LIBRARY SEARCH (LIBRARYNB)
09/29/94 19:28:00 + 38:51
SAMPLE: 4760-001-02 1644.1 09/21/94
CONDS.: EPA METHOD 8240
ENHANCED (S 15B 2N 0T)

DATA: C8696 #1542
CALI: C8696 # 3

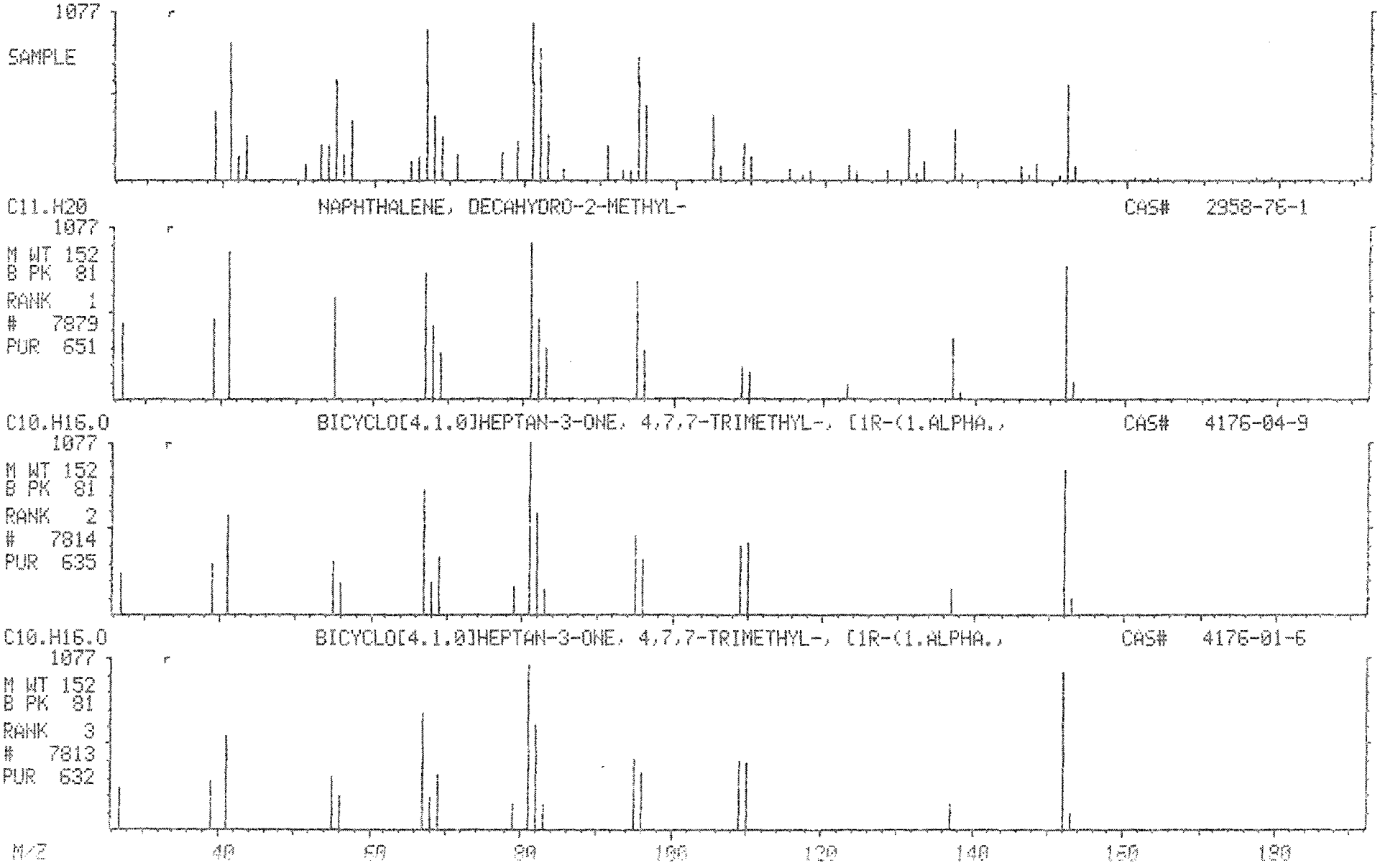
BASE M/Z: 83
RIC: 54080.



MID LIBRARY SEARCH (LIBRARYNB)
09/29/94 19:28:00 + 39:24
SAMPLE: 4760-001-02 1644.1 09/21/94
CONDS.: EPA METHOD 8240
ENHANCED (S 15B 2N 0T)

DATA: C8696 #1564
CALI: C8696 # 3

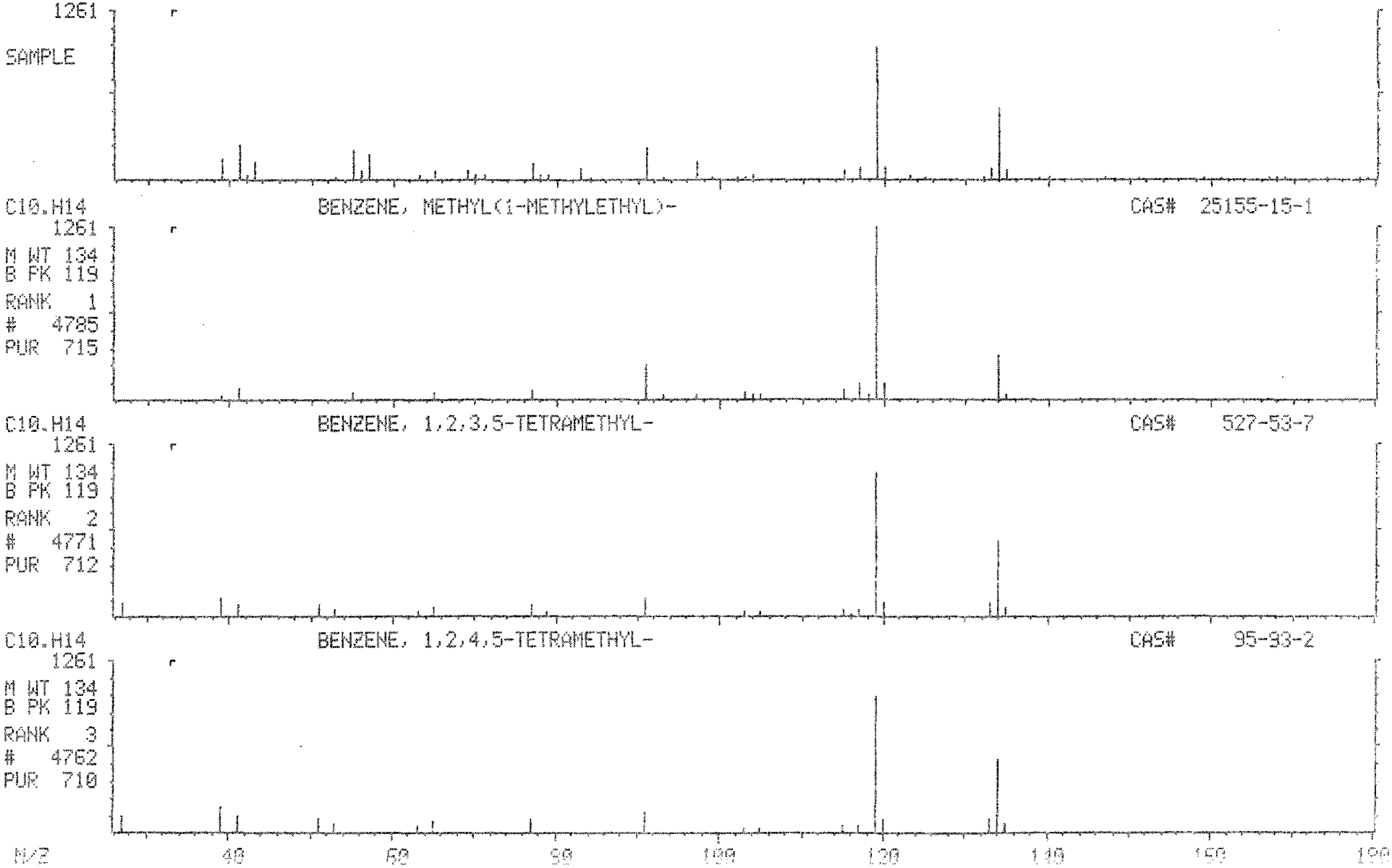
BASE M/Z: 81
RIC: 58048.



MID LIBRARY SEARCH (LIBRARYNB)
09/29/94 19:28:00 + 39:37
SAMPLE: 4760-001-02 1644.1 09/21/94
CONDS.: EPA METHOD 8240
ENHANCED (S 15B 2N 0T)

DATA: C8696 #1573
CALI: C8696 # 3

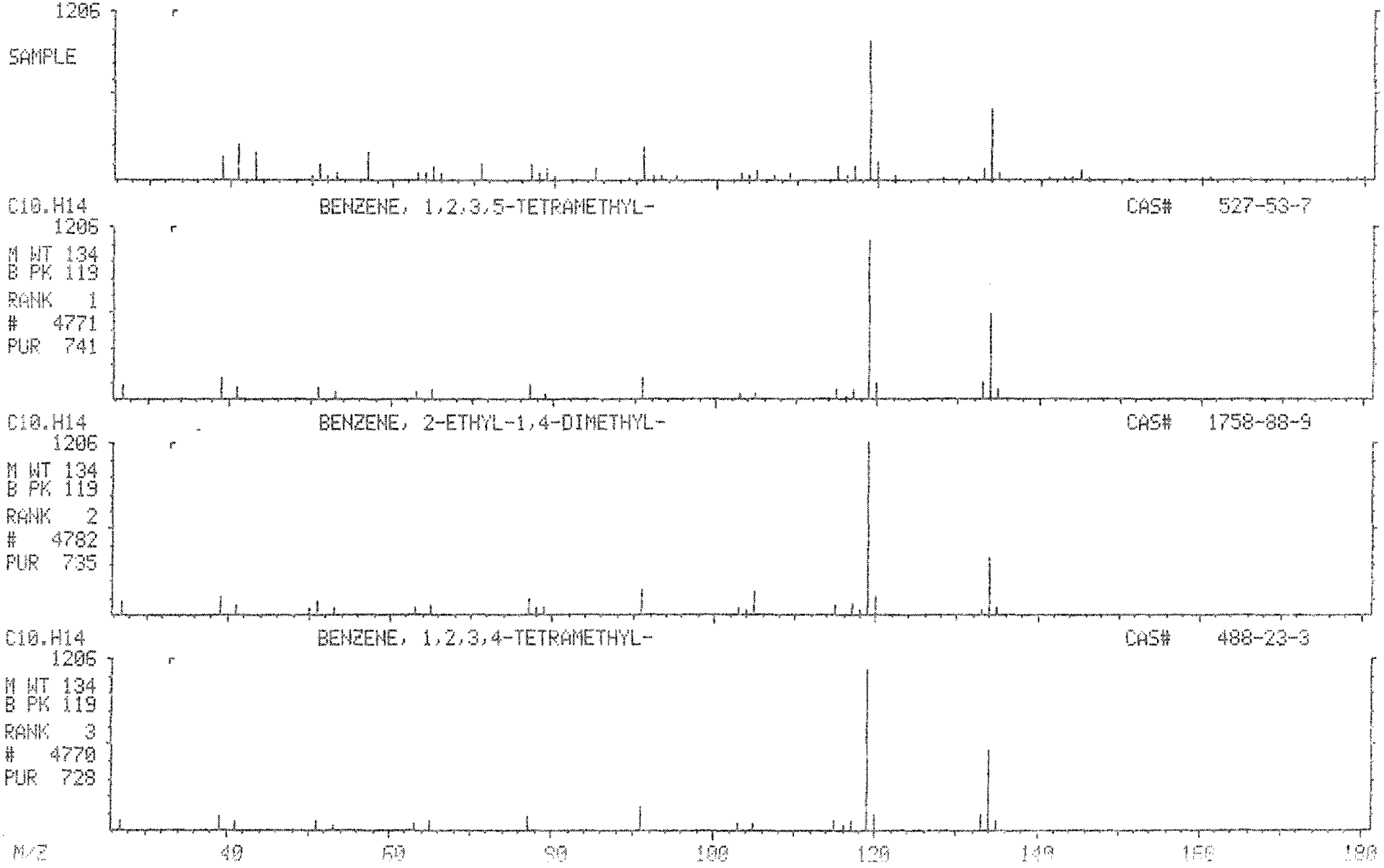
BASE M/Z: 119
RIC: 24704.



MID LIBRARY SEARCH (LIBRARYNB)
09/29/94 19:28:00 + 39:56
SAMPLE: 4760-001-02 1644.1 09/21/94
CONDS.: EPA METHOD 8240
ENHANCED (S 15B 2N 0T)

DATA: C8696 #1585
CALI: C8696 # 3

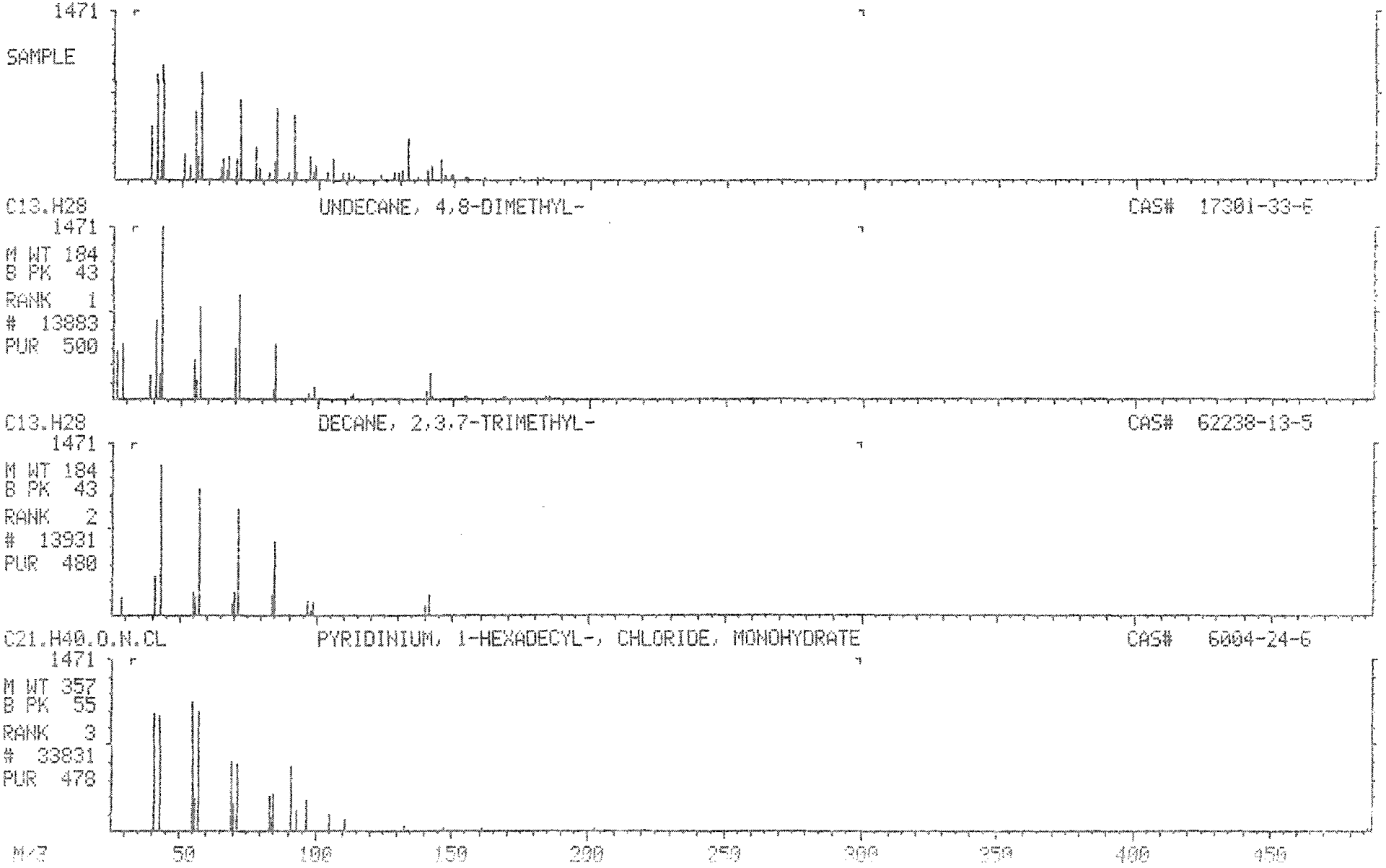
BASE M/Z: 119
RIC: 39488.

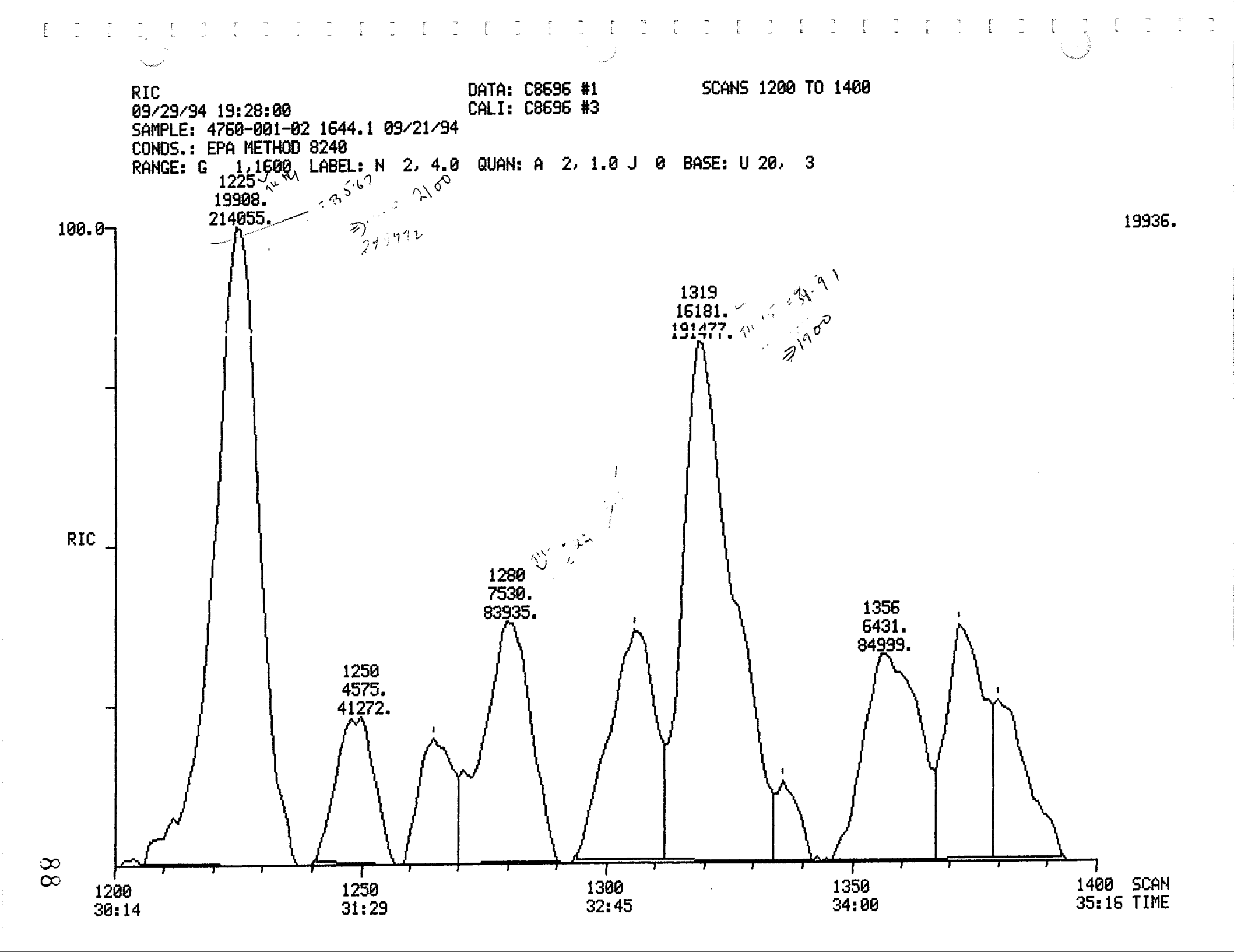


MID LIBRARY SEARCH (LIBRARYNB)
09/29/94 19:28:00 + 39:59
SAMPLE: 4760-001-02 1644.1 09/21/94
CONDS.: EPA METHOD 8240
ENHANCED (S 15B 2N 0T)

DATA: C8696 #1587
CALI: C8696 # 3

BASE M/Z: 43
RIC: 22176.




RIC
09/29/94 19:28:00
SAMPLE: 4760-001-02 1644.1 09/21/94
CONDS.: EPA METHOD 8240
RANGE: G 1,1600 LABEL: N 2, 4.0 QUAN: A 2, 1.0 J 0 BASE: U 20, 3
DATA: C8696 #1
CALI: C8696 #3
SCANS 1200 TO 1400
1225
19908.
214055.
1319
16181.
191477.
1280
7530.
83935.
1250
4575.
41272.
1356
6431.
84999.
19936.
100.0
RIC
1200
30:14
1250
31:29
1300
32:45
1350
34:00
1400 SCAN
35:16 TIME

Library Search Data: C8696 #1225 Base m/z: 43
09/29/94 19:28:00 + 30:51 Cali: C8696 # 3 RIC: 15104.
Sample: 4760-001-02 1644.1 09/21/94
Conds.: EPA METHOD 8240
Enhanced (S 15B 2N OT)

42223 spectra in LIBRARYNB searched for maximum FIT
176 matched at least 7 of the 16 largest peaks in the unknown

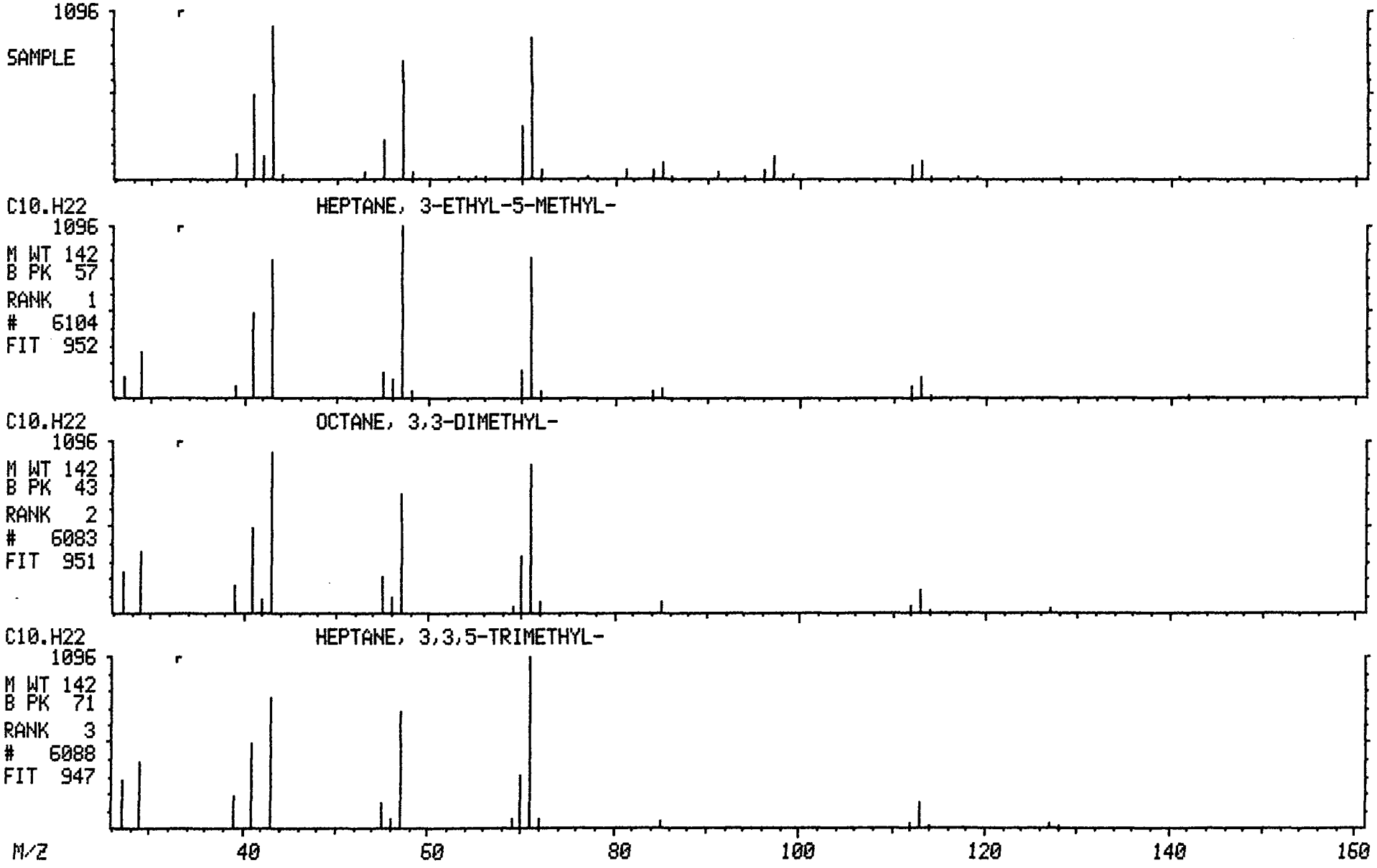
Rank	In.	Name
1	6104	HEPTANE, 3-ETHYL-5-METHYL-
2	6083	OCTANE, 3,3-DIMETHYL-
3	6088	HEPTANE, 3,3,5-TRIMETHYL-
4	6079	HEPTANE, 2,5,5-TRIMETHYL-
5	6089	HEPTANE, 5-ETHYL-2-METHYL-

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C10.H22	142	57	795	952	817
2	C10.H22	142	43	792	951	827
3	C10.H22	142	71	770	947	802
4	C10.H22	142	43	781	943	812
5	C10.H22	142	57	780	942	816

MID LIBRARY SEARCH (LIBRARYNB)
09/29/94 19:28:00 + 30:51
SAMPLE: 4760-001-02 1644.1 09/21/94
CONDS.: EPA METHOD 8240
ENHANCED (S 15B 2N 0T)

DATA: C8696 #1225
CALI: C8696 # 3

BASE M/Z: 43
RIC: 15104.



Library Search Data: C8696 #1319 Base m/z: 83
09/29/94 19:28:00 + 33:14 Cali: C8696 # 3 RIC: 12256.
Sample: 4760-001-02 1644.1 09/21/94
Conds.: EPA METHOD 8240
Enhanced (S 15B 2N OT)

42223 spectra in LIBRARYNB searched for maximum FIT
319 matched at least 7 of the 16 largest peaks in the unknown

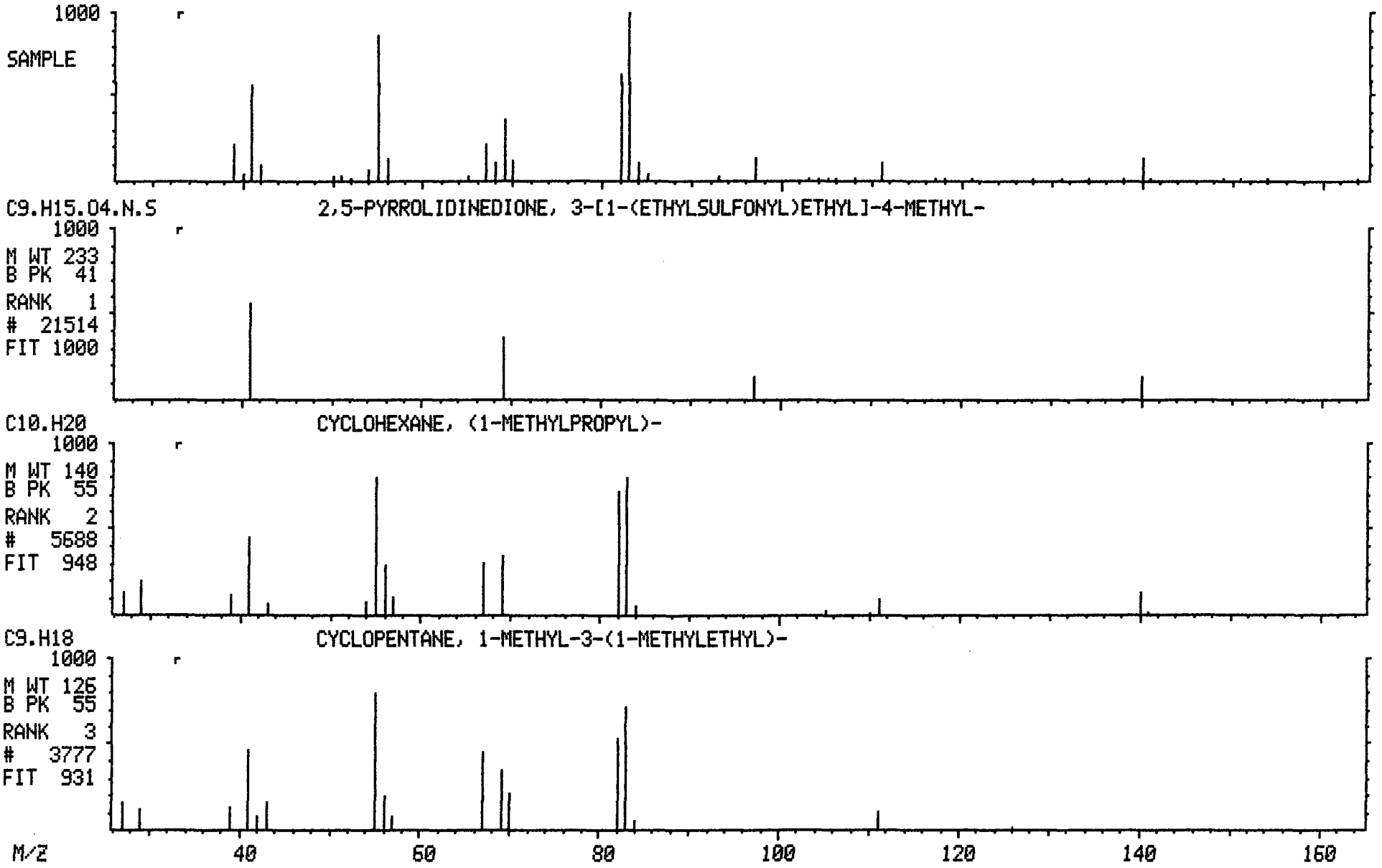
Rank	In.	Name
1	21514	2,5-PYRROLIDINEDIONE, 3-[1-(ETHYLSULFONYL)ETHYL]-4-METHYL-
2	5688	CYCLOHEXANE, (1-METHYLPROPYL)-
3	3777	CYCLOPENTANE, 1-METHYL-3-(1-METHYLETHYL)-
4	9936	CYCLOHEXANE, BROMO-
5	470	1-AZABICYCLO[3.1.0]HEXANE

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C9.H15.O4.N.S	233	41	218	1000	218
2	C10.H20	140	55	790	948	820
3	C9.H18	126	55	744	931	779
4	C6.H11.BR	162	83	502	930	526
5	C5.H9.N	83	55	600	926	608

MID LIBRARY SEARCH (LIBRARYNB)
09/29/94 19:28:00 + 33:14
SAMPLE: 4760-001-02 1644.1 09/21/94
CONDS.: EPA METHOD 8240
ENHANCED (S 15B 2N 0T)

DATA: C8696 #1319
CALI: C8696 # 3

BASE M/Z: 83
RIC: 12256.



Princeton Testing Laboratory Inc.

P.O. Box 3108
3490 U.S. Route 1
Princeton, NJ 08543-3108
(609) 452-9050
(FAX) (609) 452-1959

U.S. Army, Fort Monmouth N.J.
ATTN: SELFM-PW
Building 167
Fort Monmouth, New Jersey 07703-5108
Attention: Charles Appleby

Report Date: 11/01/94
Job Number: 9404760-001
Date Received: 09/23/94
Client Job No.: 1644/1645
Page: 1

Analysis: Volatile Organics, WW, SW-846 8240
Units: ug/liter

Parameters	*Sample I.D.: 1644.2 Fld Blk Bldgs 482/290 09/21/94*
Chloromethane	<10
Bromomethane	<10
Vinyl chloride	<10
Chloroethane	<10
Methylene chloride	<5.0
Acetone	<5.0
Carbon disulfide	<5.0
1,1-Dichloroethene	<5.0
1,1-Dichloroethane	<5.0
1,2-Dichloroethene (Total)	<5.0
Chloroform	<5.0
1,2-Dichloroethane	<5.0
2-Butanone	<5.0
1,1,1-Trichloroethane	<5.0
Carbon tetrachloride	<5.0
Bromodichloromethane	<5.0
1,1,2,2-Tetrachloroethane	<5.0
1,2-Dichloropropane	<5.0
trans-1,3-Dichloropropene	<5.0
Trichloroethene	<5.0
Dibromochloromethane	<5.0
1,1,2-Trichloroethane	<5.0
Benzene	<5.0
cis-1,3-Dichloropropene	<5.0
Bromoform	<5.0
2-Hexanone	<5.0
4-Methyl-2-Pentanone	<5.0
Tetrachloroethene	<5.0
Toluene	<5.0
Chlorobenzene	<5.0
Ethylbenzene	<5.0
Styrene	<5.0
Total Xylenes	<5.0

RECOVERY DATA	QC LIMITS	
1,2-Dichloroethane-d4 (Surrogate)	76-114%	96
Toluene-d8 (Surrogate)	88-110%	105
4-Bromofluorobenzene (Surrogate)	86-115%	92

EPA SAMPLE NO.
1644.2 09/21

1E
VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

Lab Name: Princeton Testing Lab US_ARMY,FORT_MONMOUTH.

Lab Code: PTL Case No.:4760-001 SAS No.:XXX SDG No.:XXX

Matrix: (Soil/Water) WATER Lab Sample ID: 03

Sample wt/vol: 5 (g/mL) mL Lab File ID: C8711

Level: (low/med) LOW Date Received: 09/23/94

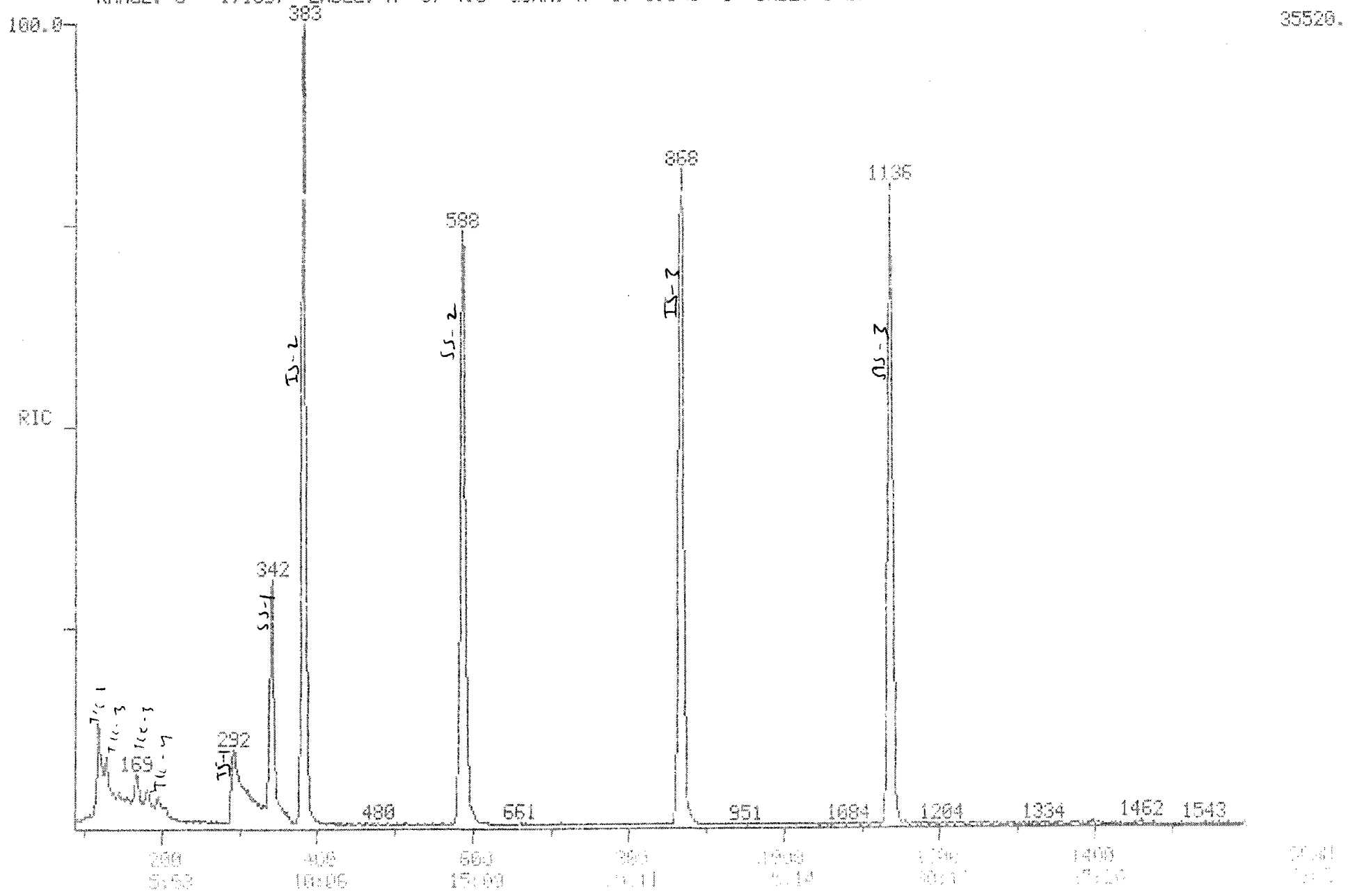
%Moisture: not dec. Date Analyzed: 10/04/94

GC Column: VOCOL ID: 0.53 mm Dilution Factor: 1

Soil Extract Vol: ul Soil Aliquot Vol: ul

Number TICs found: 4 CONCENTRATION UNITS:
(ug/L or ug/Kg) ug/L

#S	CAS NUMB	COMPOUND NAME	RT	EST. CONC.	SCAN
1	34419-766	UNKNOWN_HYDROCARBON	3:00	27	119
2	109-66-0	PENTANE_(ACN)_(DOT)	3:17	13	130
3	107-83-5	PENTANE,2-METHYL-	4:16	11	169
4	557-11-9	UREA,2-PROPENYL-	4:36	4.5	182


RIC
10/04/94 14:21:00
SAMPLE: 4760-001-03 1644.2
CONDS.: EPA METHOD 8240
RANGE: G 1,1597 LABEL: N 0, 4.0 QUAN: A 0, 1.0 J 0 BASE: U 30, 3
DATA: C8711 #1
CALI: C8711 #3
SCANS 88 TO 1597
100.0
35520.
RIC
383
588
868
1135
342
292
169
480
661
951
1084
1204
1334
1462
1543

Quantitation Report File: C8711

Data: C8711.TI
10/04/94 14:21:00
Sample: 4760-001-03 1644.2
Conds.: EPA METHOD 8240
Formula: 5ML Instrument: FINN Weight: 0.000
Submitted by: USARMY Analyst: UC Acct. No.: 4760-001

AMOUNT=AREA * REF AMNT/(REF AREA * RESP FACT)
Resp. fac. from Library Entry

No	Name
1	CI01 BROMOCHLOROMETHANE **INT. STD. **
2	CI10 1,4-DIFLUOROBENZENE **INT. STD. **
3	CI20 CHLOROBENZENE-D5 **INT. STD. **
4	CS15 1,2-DICHLOROETHANE-D4 **S. STD. **
5	CS05 TOLUENE-D8 **S. STD. **
6	CS10 4-BROMOFLUOROBENZENE **S. STD. **
7	C010 CHLOROMETHANE **
8	C015 BROMOMETHANE
9	C020 VINYL CHLORIDE *
10	C025 CHLOROETHANE
11	C030 METHYLENE CHLORIDE
12	C251 ACROLIN
13	C035 ACETONE
14	C252 ACRYLONITRILE
15	C040 CARBON DISULFIDE
16	C045 1,1-DICHLOROETHENE *
17	C050 1,1-DICHLOROETHANE **
18	C055 TRANS-1,2-DICHLOROETHENE
19	C000 TRICHLOROFLUOROMETHANE
20	C060 CHLOROFORM *
21	C065 1,2-DICHLOROETHANE
22	C110 2-BUTANONE
23	C115 1,1,1-TRICHLOROETHANE
24	C120 CARBON TETRACHLORIDE
25	C125 VINYL ACETATE
26	C130 BROMO DICHLOROMETHANE
27	C140 1,2-DICHLOROPROPANE *
28	C145 TRANS-1,3 DICHLOROPROPENE
29	C150 TRICHLOROETHENE
30	C155 DIBROMOCHLOROMETHANE
31	C160 1,1,2-TRICHLOROETHANE
32	C165 BENZENE
33	C143 CIS-1,3-DICHLOROPROPENE
34	C175 2-CHLOROETHYL VINYL ETHER
35	C180 BROMOFORM **
36	C220 TETRACHLOROETHENE
37	C210 2-HEXANONE
38	C205 4-METHYL 2-PENTANONE
39	C225 1,1,2,2-TETRACHLOROETHANE **
40	C230 TOLUENE *
41	C235 CHLOROBENZENE **
42	C240 ETHYL BENZENE *
43	C245 STYRENE
44	C250 M+P-XYLENES
45	C253 1,3-DICHLOROBENZENE
46	C254 1,4-DICHLOROBENZENE
47	C255 1,2-DICHLOROBENZENE

No	Name
48	C250 O-XYLENE
49	BENZENE-D6 **S. STD. **

No	m/z	Scan	Time	Ref	RRT	Meth	Area(Hght)	Amount	%Tot
1	49	292	7:22	1	1.000	M XX	11875.	50.000 NG	12.50
2	114	383	9:40	2	1.000	M XX	87995.	50.000 NG	12.50
3	117	867	21:53	3	1.000	A BB	47327.	50.000 NG	12.50
4	65	342	8:38	1	1.171	M XX	20700.	47.860 NG	11.97
5	98	588	14:50	3	0.678	A BB	60957.	52.312 NG	13.08
6	95	1136	28:40	3	1.310	A BB	42206.	46.126 NG	11.53
7	NOT FOUND								
8	NOT FOUND								
9	NOT FOUND								
10	NOT FOUND								
11	NOT FOUND								
12	NOT FOUND								
13	NOT FOUND								
14	NOT FOUND								
15	NOT FOUND								
16	NOT FOUND								
17	63	383	9:40	1	1.312	A BB	17758.	103.738 NG	25.94
18	NOT FOUND								
19	NOT FOUND								
20	NOT FOUND								
21	NOT FOUND								
22	NOT FOUND								
23	NOT FOUND								
24	NOT FOUND								
25	NOT FOUND								
26	NOT FOUND								
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28	NOT FOUND								
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30	NOT FOUND								
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35	NOT FOUND								
36	NOT FOUND								
37	NOT FOUND								
38	NOT FOUND								
39	NOT FOUND								
40	NOT FOUND								
41	NOT FOUND								
42	NOT FOUND								
43	NOT FOUND								
44	NOT FOUND								
45	NOT FOUND								
46	NOT FOUND								
47	NOT FOUND								
48	NOT FOUND								
49	NOT FOUND								

96
105
92

PROCEDURE: FILTER/TIC DIAGNOSTIC REPORT 10/04/94 15:24:50

DATA FILE: C8711

FILTER SCAN PARAMETERS

MAX. NUMBER TICS:	15
ID TABLE ENTRIES:	539
SCAN TOLERANCE :	2
MIN RIC HT. (%):	10
FIRST SCAN :	1
LAST SCAN :	1600
TIC THRESHOLD :	600

METHOD LIBRARY & LISTS

TIC I.S. LIBRARY:	LIBRARYLS
NBS SEARCH PROC :	SERLIB
PEAK FINDER PROC:	VOME
TCA I.S. LL :	LS
FILE NAME LIST :	TCAREF2

TARGET COMPOUND ANALYSIS:

TARGETS (QUAN LIST)	IS PEAKS	TOTAL TARGET PEAKS
4	3	7

FILTER PROCESSING:

<-----REJECT PEAKS----->							
TOTAL PEAKS	< 1ST SCAN	> LAST SCAN	< MIN RIC HT	< SCAN TOL	> MAX # PEAKS	TOTAL REJECTS	TOTAL TICS
10	0	0	0	6	0	6	4

TIC PROCESSING:

	SCAN#	PURITY	FIT	MW	COMPOUND NAME (BEFORE TIC THRESHOLD)
1	119	647	778	116	1-PROPANAMINE, N,2-DIMETHYL-N-
2	130	907	912	72	PENTANE (ACN)(DDT)
3	169	886	886	86	PENTANE, 2-METHYL-
4	182	689	756	100	UREA, 2-PROPENYL-

Quantitation Report File: C0711

Data: C0711.11
4/94 14:21:00
Sample: 4760-001-03 1644.2
Misc: EPA METHOD 8240
Method: 5ML Instrument: FINN Weight: 0.000
Submitted by: USARMY Analyst: UC Acct. No.: 4760-001

AMNT=AREA * REF AMNT/(REF AREA * RESP FACT)
Resp. fac. from Library Entry

No	CAS #	Name
1	0-00-0	C101 BROMOCHLOROMETHANE **INT. STD. **
2	0-00-0	C110 1,4-DIFLUOROBENZENE **INT. STD. **
3	0-00-0	C120 CHLOROBENZENE-D5 **INT. STD. **
4	34419-76-6	1-PROPANAMINE, N,2-DIMETHYL-N-NITROSO-
5	109-66-0	PENTANE (ACN)(DOT)
6	107-83-5	PENTANE, 2-METHYL-
7	557-11-9	UREA, 2-PROPENYL-

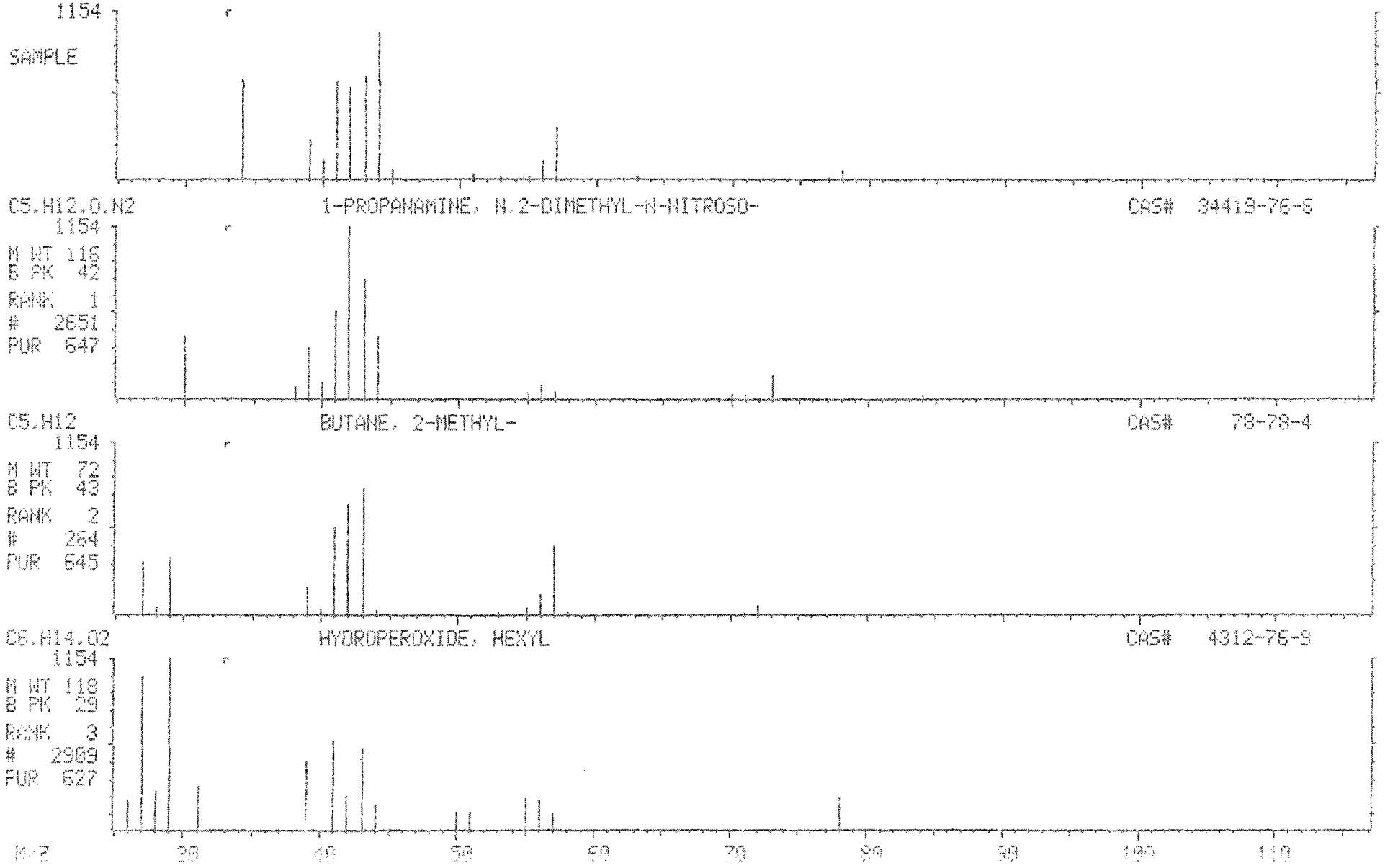
No	m/z	Scan	Time	Ref	RRT	Meth	Area(Hght)	Amount		%Tot
1	TOT	292	7:22	0	ISINV	A BB	29649.	******	UG/L	00.00
2	TOT	383	9:40	0	ISINV	A BB	183960.	******	UG/L	00.00
3	TOT	867	21:53	0	ISINV	A BB	157797.	******	UG/L	00.00
4	TOT	119	3:00	1	0.408	A BV	15753.	26.565		48.81
5	TOT	130	3:17	1	0.445	A VB	7476.	12.608		23.16
6	TOT	169	4:16	1	0.579	A BB	6390.	10.776		19.80
7	TOT	182	4:36	1	0.623	A BB	2657.	4.481		8.23

No	Ret(L)	Ratio	RRT(L)	Ratio	Amnt	Amnt(L)	R.Fac	R.Fac(L)	Ratio
1	9:13	0.80	1.000						
2	19:21	0.50	1.000						
3	23:54	0.92	1.000						
1					26.57	1.00	26.565	1.000	26.57
2					12.61	1.00	12.608	1.000	12.61
3					10.78	1.00	10.776	1.000	10.78
4					4.48	1.00	4.481	1.000	4.48

MID LIBRARY SEARCH (LIBRARYNB)
10/04/94 14:21:00 + 3:00
SAMPLE: 4760-001-03 1644.2
CONDS.: EPA METHOD 8240
ENHANCED (S 15B 2N 0T)

DATA: C8711 # 119
CALI: C8711 # 3

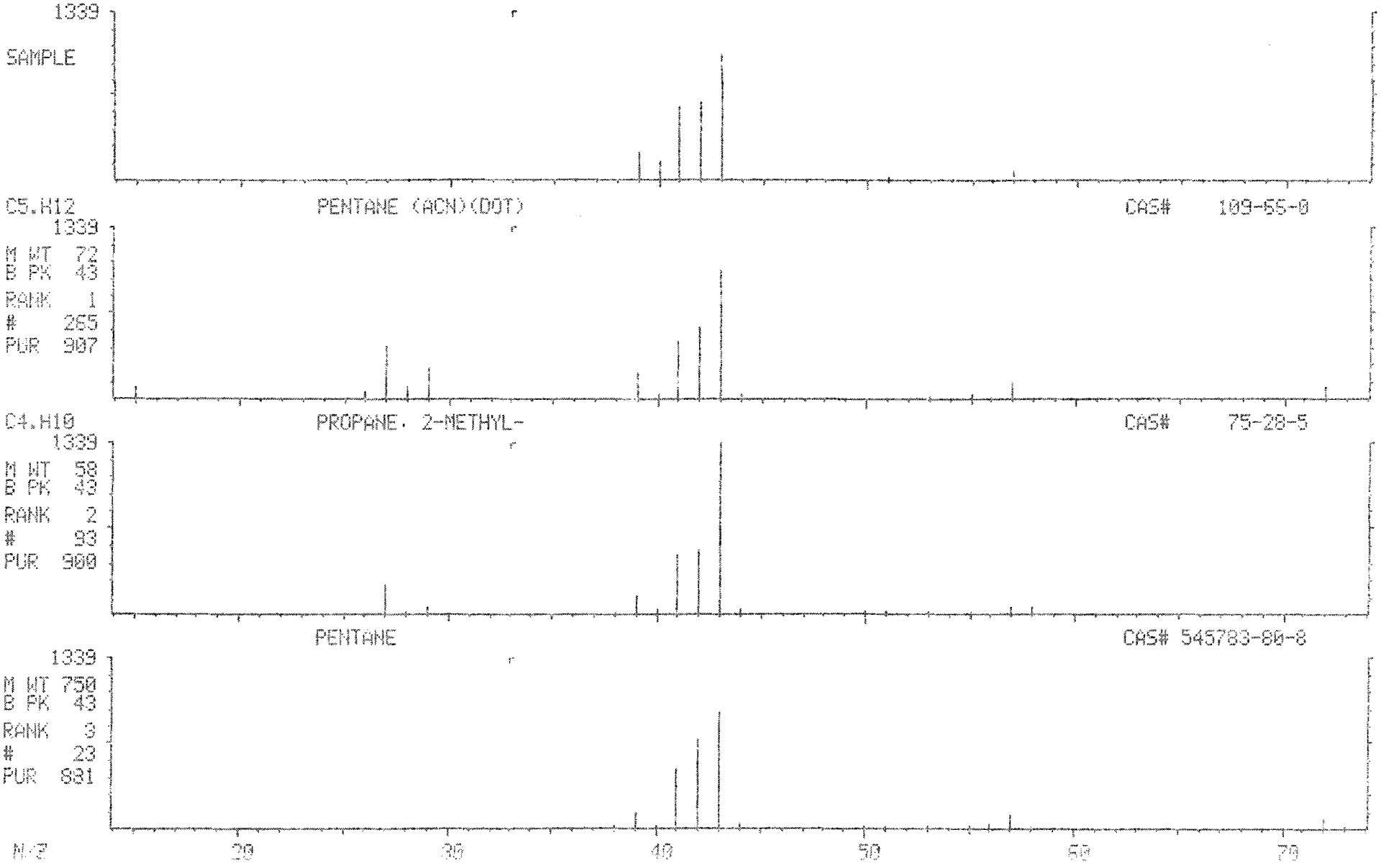
BASE M/Z: 44
RIC: 3136.



MID LIBRARY SEARCH (LIBRARYNB)
10/04/94 14:21:00 + 3:17
SAMPLE: 4760-001-03 1644.2
CONDS.: EPA METHOD 8240
ENHANCED (S 15B 2N 0T)

DATA: C8711 # 130
CALI: C8711 # 3

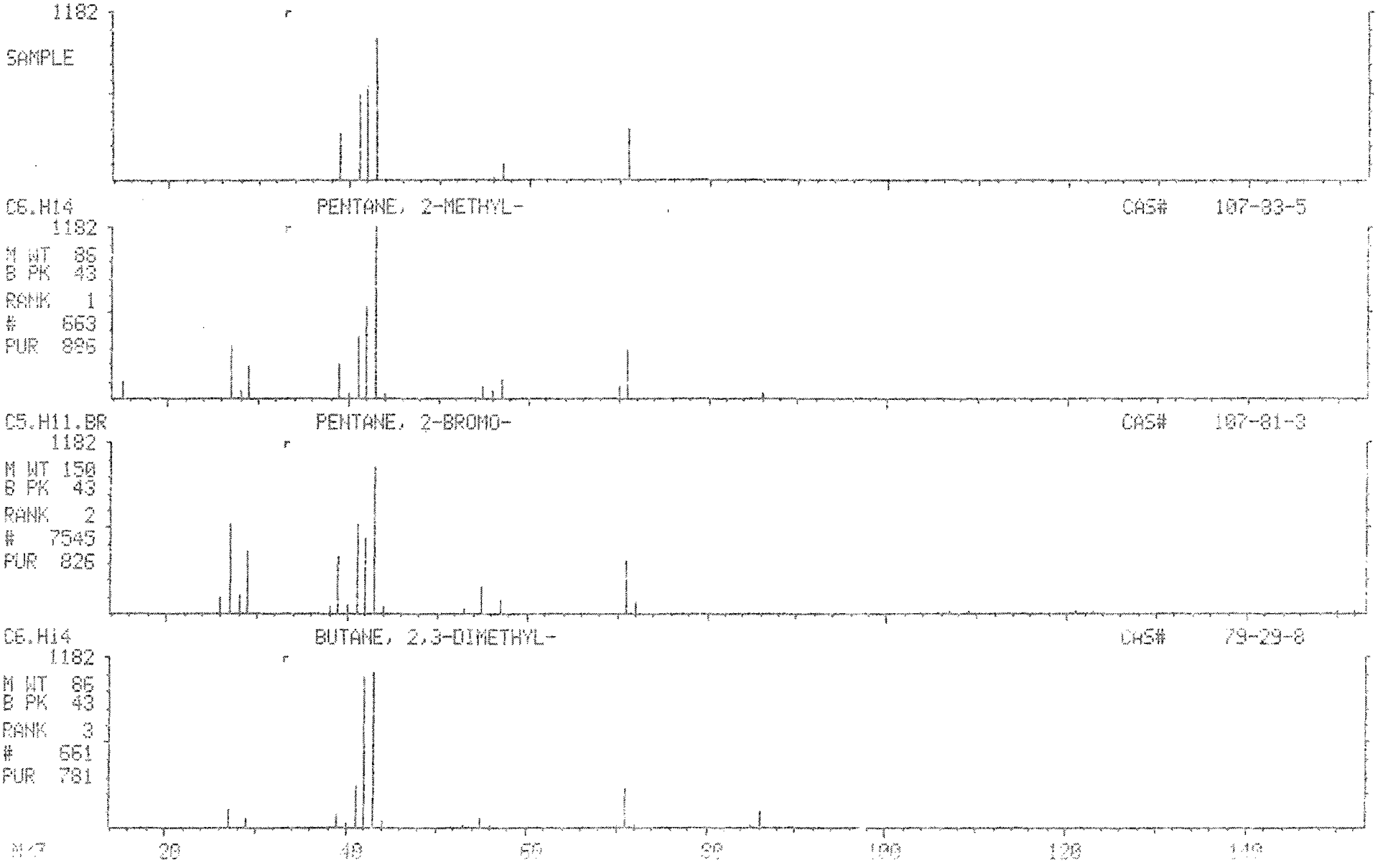
BASE M/Z: 43
RIC: 1262.



MID LIBRARY SEARCH (LIBRARYNB)
10/04/94 14:21:00 + 4:16
SAMPLE: 4760-001-03 1644.2
CONDS.: EPA METHOD 8240
ENHANCED (S 15B 2N 0T)

DATA: C8711 # 169
CALI: C8711 # 3

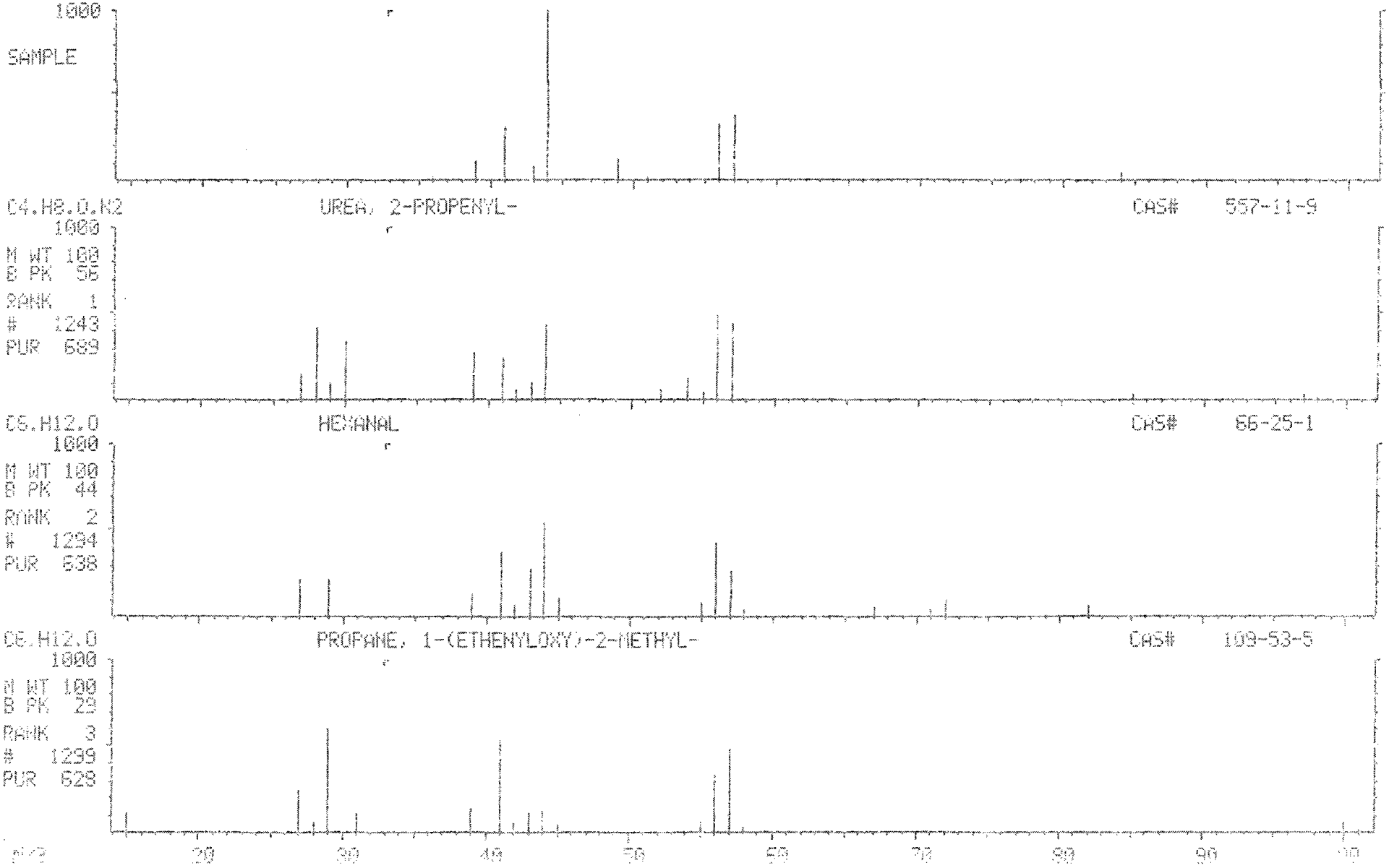
BASE M/Z: 43
RIC: 1044.



MID LIBRARY SEARCH (LIBRARYNB)
10/04/94 14:21:00 + 4:36
SAMPLE: 4760-001-03 1644.2
CONDS.: EPA METHOD 8240
ENHANCED (S 15B 2N 0T)

DATA: C8711 # 182
CALI: C8711 # 3

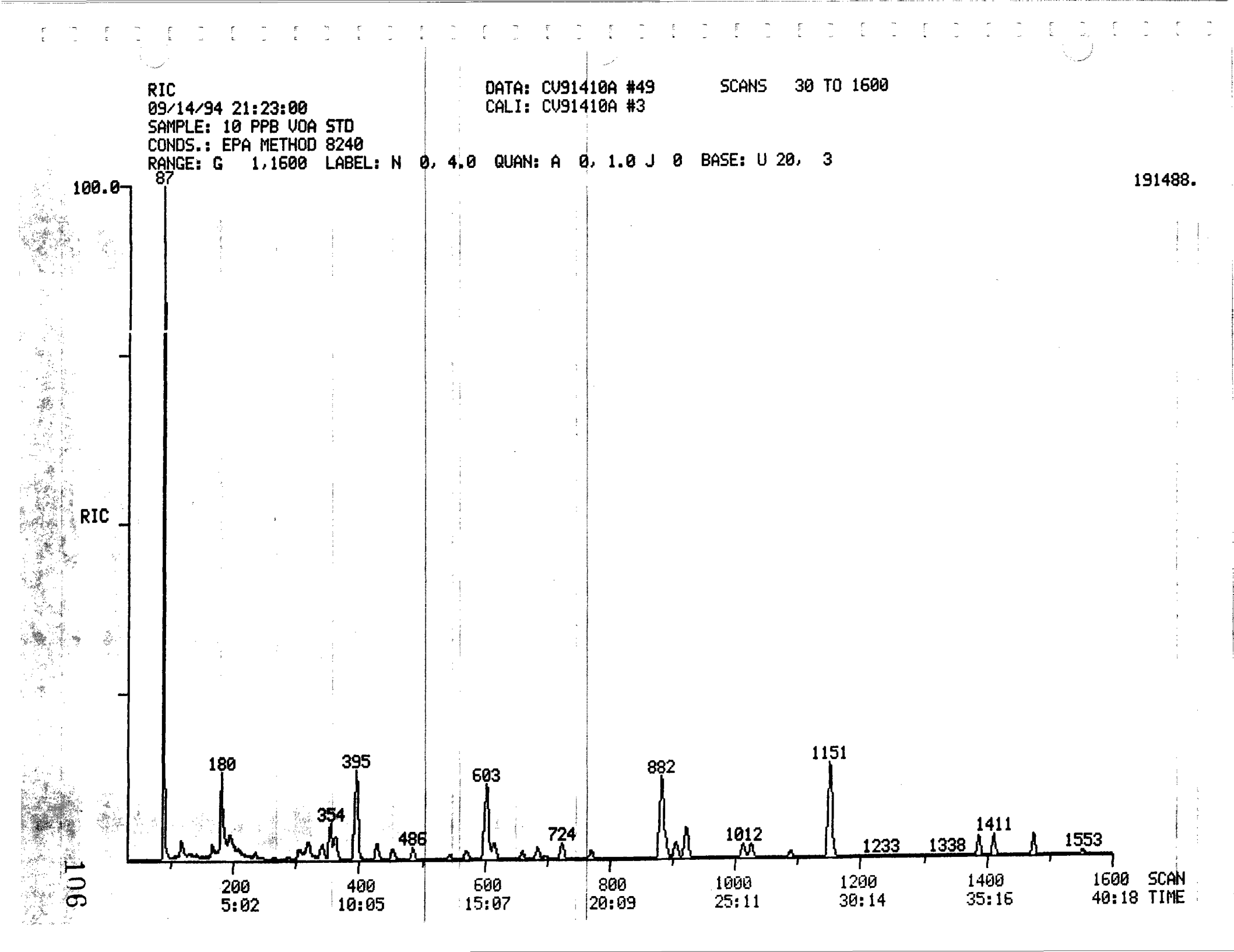
BASE M/Z: 44
RIC: 963.



Initial Calibration Data

Instrument Identifier: FINN
Calibration Date: 09/14/94
%5% RSD - 8240.
MIN RF FOR SPCC (**) = 0.30

	Compound	CV 9/14/10A RF 10	CV 9/14 20A RF 20	CV 9/14 50A RF 50	CV 9/14 100 RF 100	CV 9/14 200 RF 200	Mean RF	%RSD
C010	CHLOROMETHANE **	1.686	2.462	2.506	2.683	2.223	2.312	16.727
C015	BROMOMETHANE	2.211	2.079	2.357	3.058	2.719	2.485	16.092
C020	VINYL CHLORIDE *	2.611	2.681	2.617	2.465	1.693	2.413	17.003
C025	CHLOROETHANE	0.714	0.669	0.817	1.245	1.155	0.920	28.576
C030	METHYLENE CHLORIDE	4.443	3.771	3.290	3.207	3.887	3.720	13.455
C251	ACROLIN	0.346	0.139	0.081	0.119	0.267	0.190	58.625
C035	ACETONE	1.737	1.058	1.339	1.617	1.333	1.417	18.830
C252	ACRYLONITRILE	0.342	0.518	0.525	0.612	0.568	0.513	20.021
C040	CARBON DISULFIDE	7.147	7.850	7.536	8.086	9.395	8.003	10.673
C045	1,1-DICHLOROETHENE *	0.706	0.553	0.594	0.720	0.815	0.678	15.527
C050	1,1-DICHLOROETHANE **	1.042	0.684	0.721	0.736	0.851	0.807	18.036
C055	TRANS-1,2-DICHLOROETHEN	2.335	2.026	2.153	2.326	2.518	2.272	8.296
C000	TRICHLOROFLUOROMETHA	0.722	0.511	0.524	0.524	0.724	0.601	18.590
C060	CHLOROFORM *	1.605	0.970	0.945	0.935	1.164	1.124	25.345
C065	1,2-DICHLOROETHANE	3.583	2.634	2.330	2.424	2.650	2.724	18.322
C110	2-BUTANONE	0.185	0.109	0.108	0.119	0.144	0.133	24.410
C115	1,1,1-TRICHLOROETHANE	0.674	0.680	0.712	0.728	0.759	0.711	4.965
C120	CARBON TETRACHLORIDE	0.541	0.544	0.545	0.571	0.604	0.561	4.811
C125	VINYL ACETATE	0.343	0.279	0.504	0.618	0.586	0.466	32.005
C130	BROMO DICHLOROMETHANE	0.583	0.634	0.697	0.697	0.727	0.668	8.680
140	1,2-DICHLOROPROPANE	0.322	0.350	0.357	0.363	0.368	0.352	5.194
C145	TRANS-1,3 DICHLOROPROPE	0.407	0.483	0.539	0.552	0.568	0.510	12.924
C150	TRICHLOROETHENE	0.418	0.433	0.460	0.439	0.451	0.440	3.707
C155	DIBROMOCHLOROMETHANE	0.437	0.447	0.497	0.509	0.522	0.482	7.922
C160	1,1,2-TRICHLOROETHANE	0.315	0.319	0.336	0.331	0.342	0.329	3.427
C165	BENZENE	0.924	0.984	1.010	1.041	1.014	0.994	4.447
C143	CIS-1,3-DICHLOROPROPENE	0.377	0.429	0.486	0.501	0.497	0.458	11.752
C175	2-CHLOROETHYL VINYL ETH	0.000	0.000	0.180	0.194	0.204	0.193	6.26
C180	BROMOFORM **	0.271	0.293	0.303	0.308	0.356	0.306	10.221
C220	TETRACHLOROETHENE	0.507	0.489	0.517	0.490	0.500	0.501	2.380
C210	2-HEXANONE	0.419	0.508	0.497	0.528	0.291	0.449	21.730
C205	4-METHYL 2-PENTANONE	0.320	0.226	0.223	0.247	0.229	0.249	16.299
C225	1,1,2,2-TETRACHLOROETHA	0.617	0.632	0.606	0.504	0.621	0.596	8.764
C230	TOLUENE *	1.375	1.326	1.446	1.405	1.408	1.392	3.215
C235	CHLOROBENZENE **	0.957	0.940	0.985	0.984	0.976	0.968	2.018
C240	ETHYL BENZENE *	0.388	0.409	0.426	0.446	0.452	0.424	6.268
C245	STYRENE	0.812	0.845	0.942	0.990	1.035	0.925	10.206
C250	M+P-XYLENES	0.988	1.058	1.138	1.199	1.127	1.102	7.337
C253	1,3-DICHLOROBENZENE	0.867	0.916	0.947	0.860	0.905	0.899	4.027
C254	1,4-DICHLOROBENZENE	0.888	0.927	0.961	0.857	0.915	0.909	4.330
C255	1,2-DICHLOROBENZENE	0.832	0.921	0.880	0.770	0.849	0.850	6.617
C250	O-XYLENE	0.512	0.572	0.557	0.595	0.558	0.559	5.462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09/14/94 21:23:00
SAMPLE: 10 PPB VOA STD
CONDS.: EPA METHOD 8240
RANGE: G 1,1600 LABEL: N 0, 4.0 QUAN: A 0, 1.0 J 0 BASE: U 20, 3
DATA: CV91410A #49
CALI: CV91410A #3
SCANS 30 TO 1600
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100.0
87
RIC
180
354
395
486
603
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400
10:05
600
15:07
800
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1000
25:11
1200
30:14
1400
35:16
1600
40:18
SCAN
TIME
106

Quantitation Report File: CV91410A

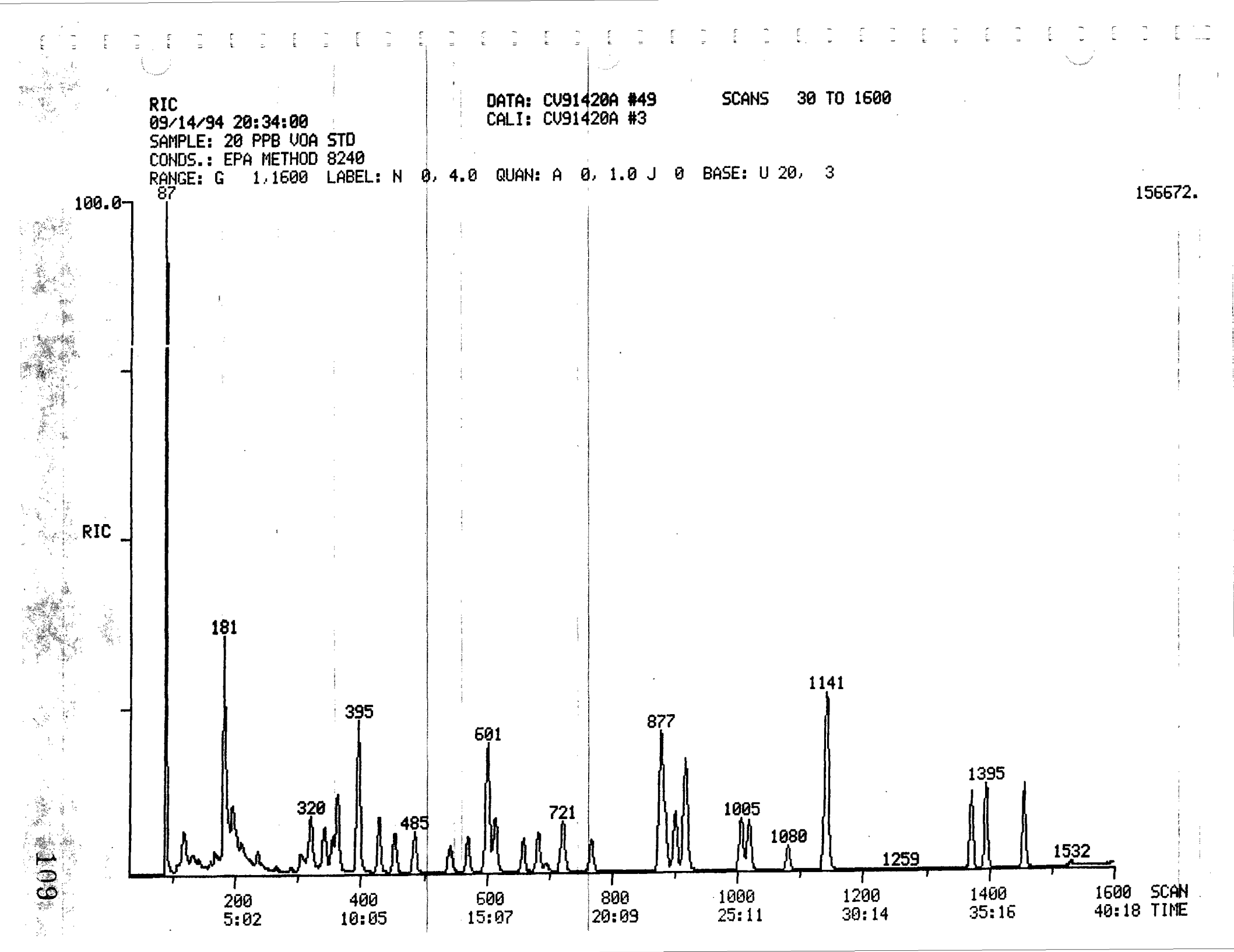
Data: CV91410A.TI
9/14/94 21:23:00
Sample: 10 PPB VOA STD
Conds.: EPA METHOD 8240
Formula: 5ML Instrument: FINN Weight: 0.000
Submitted by: PTL Analyst: UC Acct. No.: 5-PT

AMOUNT=AREA * REF AMNT/(REF AREA * RESP FACT)
Resp. fac. from Library Entry

No	Name
1	CI01 BROMOCHLOROMETHANE **INT. STD. **
2	CI10 1,4-DIFLUOROBENZENE **INT. STD. **
3	CI20 CHLOROBENZENE-D5 **INT. STD. **
4	CS15 1,2-DICHLOROETHANE-D4 **S. STD. **
5	CS05 TOLUENE-D8 **S. STD. **
6	CS10 4-BROMOFLUOROBENZENE **S. STD. **
7	C010 CHLOROMETHANE **
8	C015 BROMOMETHANE
9	C020 VINYL CHLORIDE *
10	C025 CHLOROETHANE
11	C030 METHYLENE CHLORIDE
12	C251 ACROLIN
13	C035 ACETONE
14	C252 ACRYLONITRILE
15	C040 CARBON DISULFIDE
16	C045 1,1-DICHLOROETHENE *
17	C050 1,1-DICHLOROETHANE **
18	C055 TRANS-1,2-DICHLOROETHENE
19	C000 TRICHLOROFLUOROMETHANE
20	C060 CHLOROFORM *
21	C065 1,2-DICHLOROETHANE
22	C110 2-BUTANONE
23	C115 1,1,1-TRICHLOROETHANE
24	C120 CARBON TETRACHLORIDE
25	C125 VINYL ACETATE
26	C130 BROMO DICHLOROMETHANE
27	C140 1,2-DICHLOROPROPANE *
28	C145 TRANS-1,3 DICHLOROPROPENE
29	C150 TRICHLOROETHENE
30	C155 DIBROMOCHLOROMETHANE
31	C160 1,1,2-TRICHLOROETHANE
32	C165 BENZENE
33	C143 CIS-1,3-DICHLOROPROPENE
34	C175 2-CHLOROETHYL VINYL ETHER
35	C180 BROMOFORM **
36	C220 TETRACHLOROETHENE
37	C210 2-HEXANONE
38	C205 4-METHYL 2-PENTANONE
39	C225 1,1,2,2-TETRACHLOROETHANE **
40	C230 TOLUENE *
41	C235 CHLOROBENZENE **
42	C240 ETHYL BENZENE *
43	C245 STYRENE
44	C250 M+P-XYLENES
45	C253 1,3-DICHLOROBENZENE
46	C254 1,4-DICHLOROBENZENE
47	C255 1,2-DICHLOROBENZENE

No	Name
48	C250 O-XYLENE
49	BENZENE-D6 **S. STD. **

No	m/z	Scan	Time	Ref	RRT	Meth	Area(Hght)	Amount	%Tot
1	49	305	7:41	1	1.000	A BB	6750.	50.000 NG	6.46
2	114	395	9:57	2	1.000	A BB	64012.	50.000 NG	6.46
3	117	882	22:13	3	1.000	A BB	49854.	50.000 NG	6.46
4	65	354	8:55	1	1.161	A BB	19365.	78.771 NG	10.17
5	98	603	15:11	3	0.684	A BB	61090.	49.768 NG	6.43
6	95	1151	29:00	3	1.305	A BB	43458.	45.086 NG	5.82
7	50	107	2:42	1	0.351	A BV	2275.	6.725 NG	0.87
8	94	130	3:16	1	0.426	M XX	2984.	9.381 NG	1.21
9	62	114	2:52	1	0.374	A BB	3524.	9.975 NG	1.29
10	64	136	3:26	1	0.446	M XX	964.	8.742 NG	1.13
11	49	194	4:53	1	0.636	A BB	5997.	13.505 NG	1.74
12	56	164	4:08	1	0.538	M XX	466.	43.114 NG	5.57
13	43	166	4:11	1	0.544	A BB	2344.	12.977 NG	1.68
14	53	203	5:07	1	0.666	A BB	461.	6.516 NG	0.84
15	76	194	4:53	1	0.636	A BB	9648.	9.484 NG	1.22
16	96	169	4:15	1	0.554	M XX	953.	11.897 NG	1.54
17	63	233	5:52	1	0.764	M XX	1406.	14.447 NG	1.87
18	61	208	5:14	1	0.682	M XX	3151.	10.847 NG	1.40
19	101	142	3:35	1	0.466	A BB	974.	13.783 NG	1.78
20	83	288	7:15	1	0.944	M XX	2166.	16.991 NG	2.19
21	62	364	9:10	1	1.193	A BB	4836.	15.381 NG	1.99
22	43	266	6:42	2	0.673	A BB	2358.	17.212 NG	2.22
23	97	320	8:04	2	0.810	A BB	8617.	9.454 NG	1.22
24	117	341	8:35	2	0.863	A BB	6919.	9.920 NG	1.28
25	43	235	5:55	2	0.595	A BV	4389.	6.810 NG	0.88
26	83	486	12:15	2	1.230	A BB	7464.	8.372 NG	1.08
27	63	455	11:28	2	1.152	A BB	4110.	9.006 NG	1.16
28	75	572	14:25	2	1.448	A BB	5202.	7.542 NG	0.97
29	95	429	10:48	2	1.086	A BB	5348.	9.083 NG	1.17
30	129	770	19:24	2	1.949	M XX	5582.	8.782 NG	1.13
31	97	685	17:15	2	1.734	A BB	4027.	9.387 NG	1.21
32	78	362	9:07	2	0.916	A BB	11823.	9.146 NG	1.18
33	75	662	16:41	2	1.676	A BB	4814.	7.749 NG	1.00
34	NOT FOUND								
35	173	1089	27:26	2	2.757	A BB	3461.	8.937 NG	1.15
36	166	724	18:14	3	0.821	A BB	5054.	9.811 NG	1.27
37	43	545	13:44	3	0.618	M XX	4175.	8.426 NG	1.09
38	43	697	17:33	3	0.790	A BB	3182.	14.343 NG	1.85
39	83	1149	28:57	3	1.303	A BB	6143.	10.181 NG	1.32
40	91	615	15:30	3	0.697	A BB	13701.	9.507 NG	1.23
41	112	889	22:24	3	1.008	A BB	9538.	9.716 NG	1.25
42	106	906	22:49	3	1.027	A BB	3859.	9.092 NG	1.17
43	104	1026	25:51	3	1.163	A BB	8093.	8.624 NG	1.11
44	106	922	23:14	3	1.045	A BB	9845.	8.684 NG	1.12
45	146	1387	34:56	3	1.573	A BB	8636.	9.150 NG	1.18
46	146	1411	35:33	3	1.600	A BB	8848.	9.240 NG	1.19
47	146	1474	37:08	3	1.671	A BB	8295.	9.459 NG	1.22
48	106	1012	25:30	3	1.147	A BB	5096.	9.182 NG	1.19
49	NOT FOUND								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09/14/94 20:34:00
SAMPLE: 20 PPB VOA STD
CONDS.: EPA METHOD 8240
RANGE: G 1,1600 LABEL: N 0, 4.0 QUAN: A 0, 1.0 J 0 BASE: U 20, 3
DATA: CV91420A #49
CALI: CV91420A #3
SCANS 30 TO 1600
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100.0
RIC
87
181
320
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200
5:02
400
10:05
600
15:07
800
20:09
1000
25:11
1200
30:14
1400
35:16
1600
40:18
SCAN
TIME
109

Quantitation Report File: CV91420A

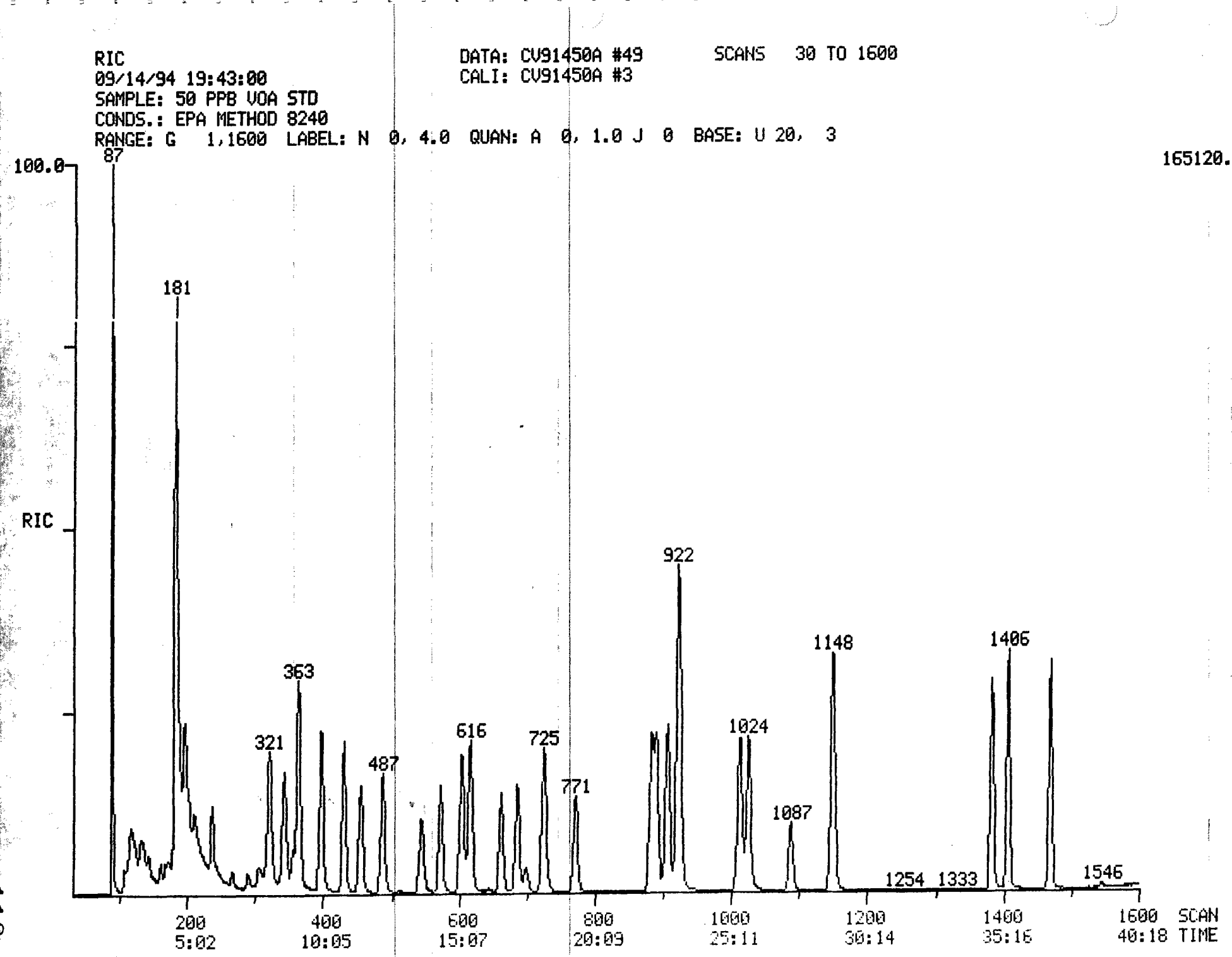
Data: CV91420A.TI
9/14/94 20:34:00
Sample: 20 PPB VOA STD
Conds.: EPA METHOD 8240
Formula: 5ML Instrument: FINN Weight: 0.000
Submitted by: PTL Analyst: UC Acct. No.: 5-PT

AMOUNT=AREA * REF AMNT/(REF AREA * RESP FACT)
Resp. fac. from Library Entry

No	Name
1	CI01 BROMOCHLOROMETHANE **INT. STD. **
2	CI10 1,4-DIFLUOROBENZENE **INT. STD. **
3	CI20 CHLOROBENZENE-D5 **INT. STD. **
4	CS15 1,2-DICHLOROETHANE-D4 **S. STD. **
5	CS05 TOLUENE-D8 **S. STD. **
6	CS10 4-BROMOFLUOROBENZENE **S. STD. **
7	C010 CHLOROMETHANE **
8	C015 BROMOMETHANE
9	C020 VINYL CHLORIDE *
10	C025 CHLOROETHANE
11	C030 METHYLENE CHLORIDE
12	C251 ACROLIN
13	C035 ACETONE
14	C252 ACRYLONITRILE
15	C040 CARBON DISULFIDE
16	C045 1,1-DICHLOROETHENE *
17	C050 1,1-DICHLOROETHANE **
18	C055 TRANS-1,2-DICHLOROETHENE
19	C000 TRICHLOROFLUOROMETHANE
20	C060 CHLOROFORM *
21	C065 1,2-DICHLOROETHANE
22	C110 2-BUTANONE
23	C115 1,1,1-TRICHLOROETHANE
24	C120 CARBON TETRACHLORIDE
25	C125 VINYL ACETATE
26	C130 BROMO DICHLOROMETHANE
27	C140 1,2-DICHLOROPROPANE *
28	C145 TRANS-1,3 DICHLOROPROPENE
29	C150 TRICHLOROETHENE
30	C155 DIBROMOCHLOROMETHANE
31	C160 1,1,2-TRICHLOROETHANE
32	C165 BENZENE
33	C143 CIS-1,3-DICHLOROPROPENE
34	C175 2-CHLOROETHYL VINYL ETHER
35	C180 BROMOFORM **
36	C220 TETRACHLOROETHENE
37	C210 2-HEXANONE
38	C205 4-METHYL 2-PENTANONE
39	C225 1,1,2,2-TETRACHLOROETHANE **
40	C230 TOLUENE *
41	C235 CHLOROBENZENE **
42	C240 ETHYL BENZENE *
43	C245 STYRENE
44	C250 M+P-XYLENES
45	C253 1,3-DICHLOROBENZENE
46	C254 1,4-DICHLOROBENZENE
47	C255 1,2-DICHLOROBENZENE

No	Name
48	C250 O-XYLENE
49	BENZENE-D6 **S. STD. **

No	m/z	Scan	Time	Ref	RRT	Meth	Area(Hght)	Amount	%Tot
1	49	305	7:41	1	1.000	A BB	8332.	50.000 NG	4.47
2	114	395	9:57	2	1.000	A BB	86467.	50.000 NG	4.47
3	117	877	22:05	3	1.000	A BB	68587.	50.000 NG	4.47
4	65	354	8:55	1	1.161	A BB	16583.	54.642 NG	4.88
5	98	601	15:08	3	0.685	A BB	84786.	50.208 NG	4.49
6	95	1141	28:45	3	1.301	A BB	60905.	45.929 NG	4.11
7	50	108	2:43	1	0.354	A BB	8204.	19.646 NG	1.76
8	94	131	3:18	1	0.430	A BB	6928.	17.643 NG	1.58
9	62	114	2:52	1	0.374	A BB	8934.	20.487 NG	1.83
10	64	134	3:23	1	0.439	A BB	2229.	16.376 NG	1.46
11	49	195	4:55	1	0.639	A BB	12567.	22.924 NG	2.05
12	56	164	4:08	1	0.538	A BV	461.	34.562 NG	3.09
13	43	167	4:12	1	0.548	A BB	3524.	15.804 NG	1.41
14	53	203	5:07	1	0.666	A BB	1726.	19.762 NG	1.77
15	76	195	4:55	1	0.639	A BB	26162.	20.832 NG	1.86
16	96	170	4:17	1	0.557	A BB	1840.	18.617 NG	1.66
17	63	234	5:54	1	0.767	A BB	2276.	18.952 NG	1.69
18	61	210	5:17	1	0.689	A BB	6752.	18.826 NG	1.68
19	101	142	3:35	1	0.466	A BB	1700.	19.487 NG	1.74
20	83	289	7:17	1	0.948	A BB	3232.	20.536 NG	1.84
21	62	363	9:09	1	1.190	A BB	8777.	22.613 NG	2.02
22	43	266	6:42	2	0.673	A BB	3756.	20.297 NG	1.81
23	97	320	8:04	2	0.810	A BB	23504.	19.089 NG	1.71
24	117	342	8:37	2	0.866	A BB	18813.	19.969 NG	1.79
25	43	235	5:55	2	0.595	A BB	9631.	11.061 NG	0.99
26	83	485	12:13	2	1.228	A BB	21901.	18.185 NG	1.63
27	63	454	11:26	2	1.149	A BB	12100.	19.626 NG	1.75
28	75	570	14:21	2	1.443	A BB	16673.	17.896 NG	1.60
29	95	428	10:47	2	1.084	A BB	14968.	18.819 NG	1.68
30	129	767	19:19	2	1.942	A BB	15438.	17.980 NG	1.61
31	97	682	17:11	2	1.727	A BB	11018.	19.013 NG	1.70
32	78	362	9:07	2	0.916	A BB	34012.	19.476 NG	1.74
33	75	659	16:36	2	1.668	A BB	14830.	17.673 NG	1.58
34	63	541	13:38	2	1.370	A BB	6204.	18.544 NG	1.66
35	173	1080	27:12	2	2.734	A BB	10122.	19.350 NG	1.73
36	166	721	18:10	3	0.822	A BB	13408.	18.919 NG	1.69
37	43	542	13:39	3	0.618	A BB	13931.	20.435 NG	1.83
38	43	695	17:30	3	0.792	A BB	6197.	20.305 NG	1.82
39	83	1140	28:43	3	1.300	A BB	17335.	20.883 NG	1.87
40	91	614	15:28	3	0.700	A BB	36363.	18.340 NG	1.64
41	112	884	22:16	3	1.008	A BB	25768.	19.079 NG	1.71
42	106	900	22:40	3	1.026	A BB	11196.	19.172 NG	1.71
43	104	1018	25:39	3	1.161	A BB	23182.	17.956 NG	1.61
44	106	916	23:04	3	1.044	A BB	29024.	18.608 NG	1.66
45	146	1372	34:34	3	1.564	A BB	25115.	19.342 NG	1.73
46	146	1395	35:08	3	1.591	A BB	25414.	19.290 NG	1.72
47	146	1456	36:41	3	1.660	A BV	25253.	20.933 NG	1.87
48	106	1005	25:19	3	1.146	A BB	15691.	20.551 NG	1.84
49	NOT FOUND								


RIC
09/14/94 19:43:00
SAMPLE: 50 PPB VOA STD
CONDS.: EPA METHOD 8240
RANGE: G 1,1600 LABEL: N 0, 4.0 QUAN: A 0, 1.0 J 0 BASE: U 20, 3
DATA: CV91450A #49
CALI: CV91450A #3
SCANS 30 TO 1600
165120.
100.0
RIC
87
181
321
363
487
616
725
771
922
1024
1087
1148
1254
1333
1406
1546
200
5:02
400
10:05
600
15:07
800
20:09
1000
25:11
1200
30:14
1400
35:16
1600
40:18
SCAN
TIME

Quantitation Report File: CV91450A

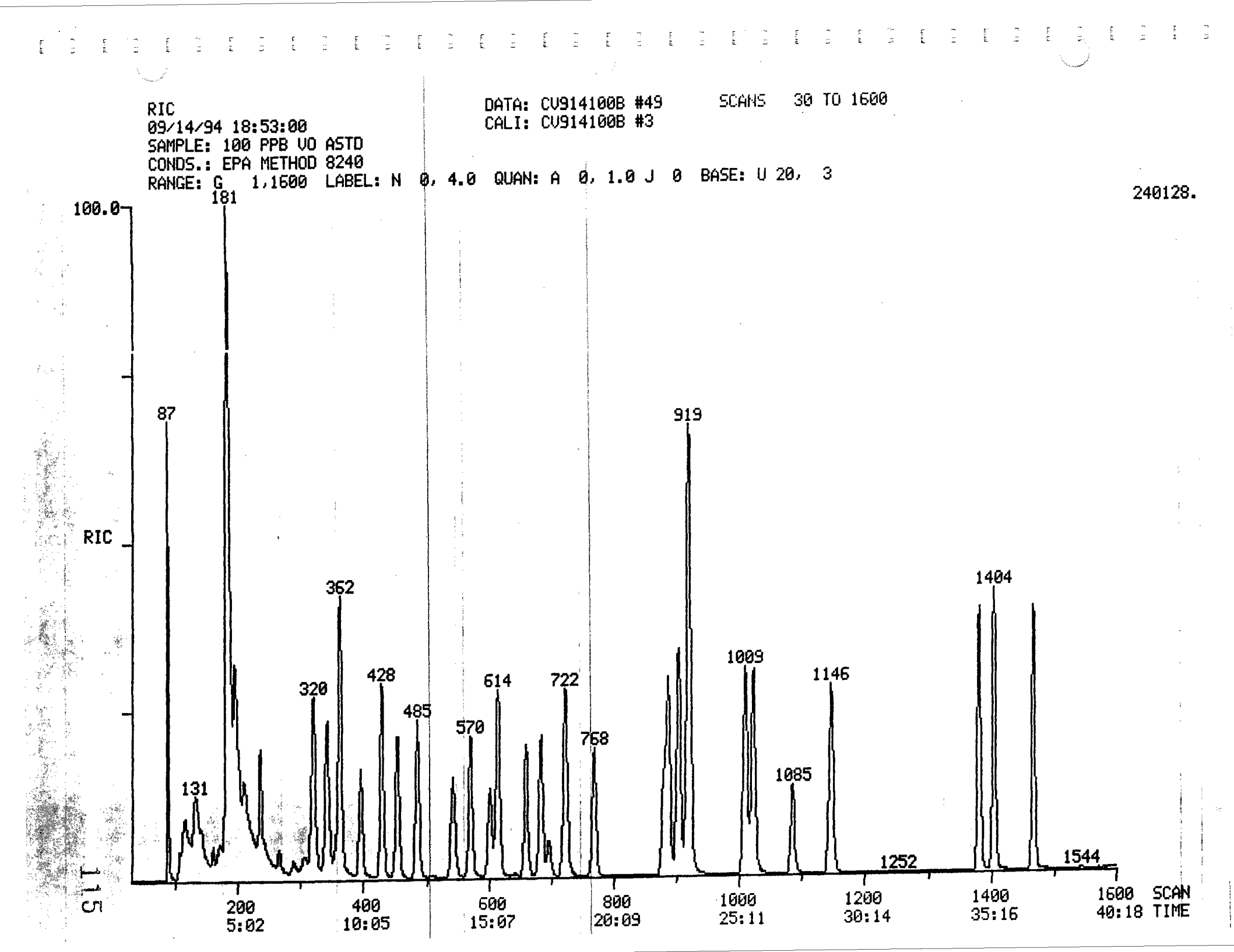
Data: CV91450A.TI
9/14/94 19:43:00
Sample: 50 PPB VOA STD
Conds.: EPA METHOD 8240
Formula: 5ML Instrument: FINN Weight: 0.000
Submitted by: PTL Analyst: UC Acct. No.: 5-PT

AMOUNT=AREA * REF AMNT/(REF AREA * RESP FACT)
Resp. fac. from Library Entry

No	Name
1	CI01 BROMOCHLOROMETHANE **INT. STD. **
2	CI10 1,4-DIFLUOROBENZENE **INT. STD. **
3	CI20 CHLOROBENZENE-D5 **INT. STD. **
4	CS15 1,2-DICHLOROETHANE-D4 **S. STD. **
5	CS05 TOLUENE-D8 **S. STD. **
6	CS10 4-BROMOFLUOROBENZENE **S. STD. **
7	C010 CHLOROMETHANE **
8	C015 BROMOMETHANE
9	C020 VINYL CHLORIDE *
10	C025 CHLOROETHANE
11	C030 METHYLENE CHLORIDE
12	C251 ACROLIN
13	C035 ACETONE
14	C252 ACRYLONITRILE
15	C040 CARBON DISULFIDE
16	C045 1,1-DICHLOROETHENE *
17	C050 1,1-DICHLOROETHANE **
18	C055 TRANS-1,2-DICHLOROETHENE
19	C000 TRICHLOROFLUOROMETHANE
20	C060 CHLOROFORM *
21	C065 1,2-DICHLOROETHANE
22	C110 2-BUTANONE
23	C115 1,1,1-TRICHLOROETHANE
24	C120 CARBON TETRACHLORIDE
25	C125 VINYL ACETATE
26	C130 BROMO DICHLOROMETHANE
27	C140 1,2-DICHLOROPROPANE *
28	C145 TRANS-1,3 DICHLOROPROPENE
29	C150 TRICHLOROETHENE
30	C155 DIBROMOCHLOROMETHANE
31	C160 1,1,2-TRICHLOROETHANE
32	C165 BENZENE
33	C143 CIS-1,3-DICHLOROPROPENE
34	C175 2-CHLOROETHYL VINYL ETHER
35	C180 BROMOFORM **
36	C220 TETRACHLOROETHENE
37	C210 2-HEXANONE
38	C205 4-METHYL 2-PENTANONE
39	C225 1,1,2,2-TETRACHLOROETHANE **
40	C230 TOLUENE *
41	C235 CHLOROBENZENE **
42	C240 ETHYL BENZENE *
43	C245 STYRENE
44	C250 M+P-XYLENES
45	C253 1,3-DICHLOROBENZENE
46	C254 1,4-DICHLOROBENZENE
47	C255 1,2-DICHLOROBENZENE

No	Name
48	C250 O-XYLENE
49	BENZENE-D6 **S. STD. **

No	m/z	Scan	Time	Ref	RRT	Meth	Area(Hght)	Amount	%Tot
1	49	305	7:41	1	1.000	A BB	9202.	50.000 NG	2.08
2	114	396	9:59	2	1.000	A BB	88266.	50.000 NG	2.08
3	117	882	22:13	3	1.000	A BB	71581.	50.000 NG	2.08
4	65	355	8:57	1	1.164	A BB	16758.	50.000 NG	2.08
5	98	603	15:11	3	0.684	A BB	88121.	50.000 NG	2.08
6	95	1149	28:57	3	1.303	A BB	69198.	50.000 NG	2.08
7	50	107	2:42	1	0.351	A BB	23060.	50.000 NG	2.08
8	94	130	3:16	1	0.426	A BB	21684.	50.000 NG	2.08
9	62	114	2:52	1	0.374	A BB	24081.	50.000 NG	2.08
10	64	134	3:23	1	0.439	A BB	7516.	50.000 NG	2.08
11	49	195	4:55	1	0.639	A BB	30271.	50.000 NG	2.08
12	56	164	4:08	1	0.538	M XX	737.	50.000 NG	2.08
13	43	166	4:11	1	0.544	A BB	12313.	50.000 NG	2.08
14	53	203	5:07	1	0.666	A BB	4823.	50.000 NG	2.08
15	76	195	4:55	1	0.639	A BB	69349.	50.000 NG	2.08
16	96	171	4:18	1	0.561	A BB	5458.	50.000 NG	2.08
17	63	235	5:55	1	0.770	A BB	6633.	50.000 NG	2.08
18	61	210	5:17	1	0.689	A BB	19804.	50.000 NG	2.08
19	101	142	3:35	1	0.466	A BB	4817.	50.000 NG	2.08
20	83	289	7:17	1	0.948	A BB	8690.	50.000 NG	2.08
21	62	365	9:12	1	1.197	A BB	21433.	50.000 NG	2.08
22	43	266	6:42	2	0.672	A BB	9445.	50.000 NG	2.08
23	97	321	8:05	2	0.811	A BB	62845.	50.000 NG	2.08
24	117	343	8:38	2	0.866	A BB	48086.	50.000 NG	2.08
25	43	235	5:55	2	0.593	A BB	44440.	50.000 NG	2.08
26	83	487	12:16	2	1.230	A BB	61468.	50.000 NG	2.08
27	63	455	11:28	2	1.149	A BB	31466.	50.000 NG	2.08
28	75	572	14:25	2	1.444	A BB	47551.	50.000 NG	2.08
29	95	429	10:48	2	1.083	A BB	40595.	50.000 NG	2.08
30	129	771	19:25	2	1.947	A BB	43823.	50.000 NG	2.08
31	97	685	17:15	2	1.730	A BB	29578.	50.000 NG	2.08
32	78	363	9:09	2	0.917	A BB	89133.	50.000 NG	2.08
33	75	662	16:41	2	1.672	A BB	42829.	50.000 NG	2.08
34	63	543	13:41	2	1.371	M XX	17076.	50.000 NG	2.08
35	173	1087	27:23	2	2.745	A BB	26699.	50.000 NG	2.08
36	166	725	18:16	3	0.822	A BB	36983.	50.000 NG	2.08
37	43	544	13:42	3	0.617	A BB	35575.	50.000 NG	2.08
38	43	698	17:35	3	0.791	A BB	15926.	50.000 NG	2.08
39	83	1148	28:55	3	1.302	A BB	43317.	50.000 NG	2.08
40	91	616	15:31	3	0.698	A BB	103464.	50.000 NG	2.08
41	112	889	22:24	3	1.008	A BB	70477.	50.000 NG	2.08
42	106	905	22:48	3	1.026	A BB	30472.	50.000 NG	2.08
43	104	1024	25:48	3	1.161	A BB	67372.	50.000 NG	2.08
44	106	922	23:14	3	1.045	A BB	81395.	50.000 NG	2.08
45	146	1382	34:49	3	1.567	A BB	67759.	50.000 NG	2.08
46	146	1406	35:25	3	1.594	A BB	68748.	50.000 NG	2.08
47	146	1468	36:59	3	1.664	A BB	62953.	50.000 NG	2.08
48	106	1011	25:28	3	1.146	A BB	39843.	50.000 NG	2.08
49	NOT FOUND								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09/14/94 18:53:00
SAMPLE: 100 PPB VO ASTD
CONDS.: EPA METHOD 8240
RANGE: G 1,1600 LABEL: N 0, 4.0 QUAN: A 0, 1.0 J 0 BASE: U 20, 3
DATA: CV914100B #49 SCANS 30 TO 1600
CALI: CV914100B #3
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100.0
RIC
87
181
131
320
362
428
485
570
614
722
768
919
1009
1085
1146
1252
1404
1544
200
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400
10:05
600
15:07
800
20:09
1000
25:11
1200
30:14
1400
35:16
1600
40:18
SCAN
TIME

Quantitation Report File: CV914100B

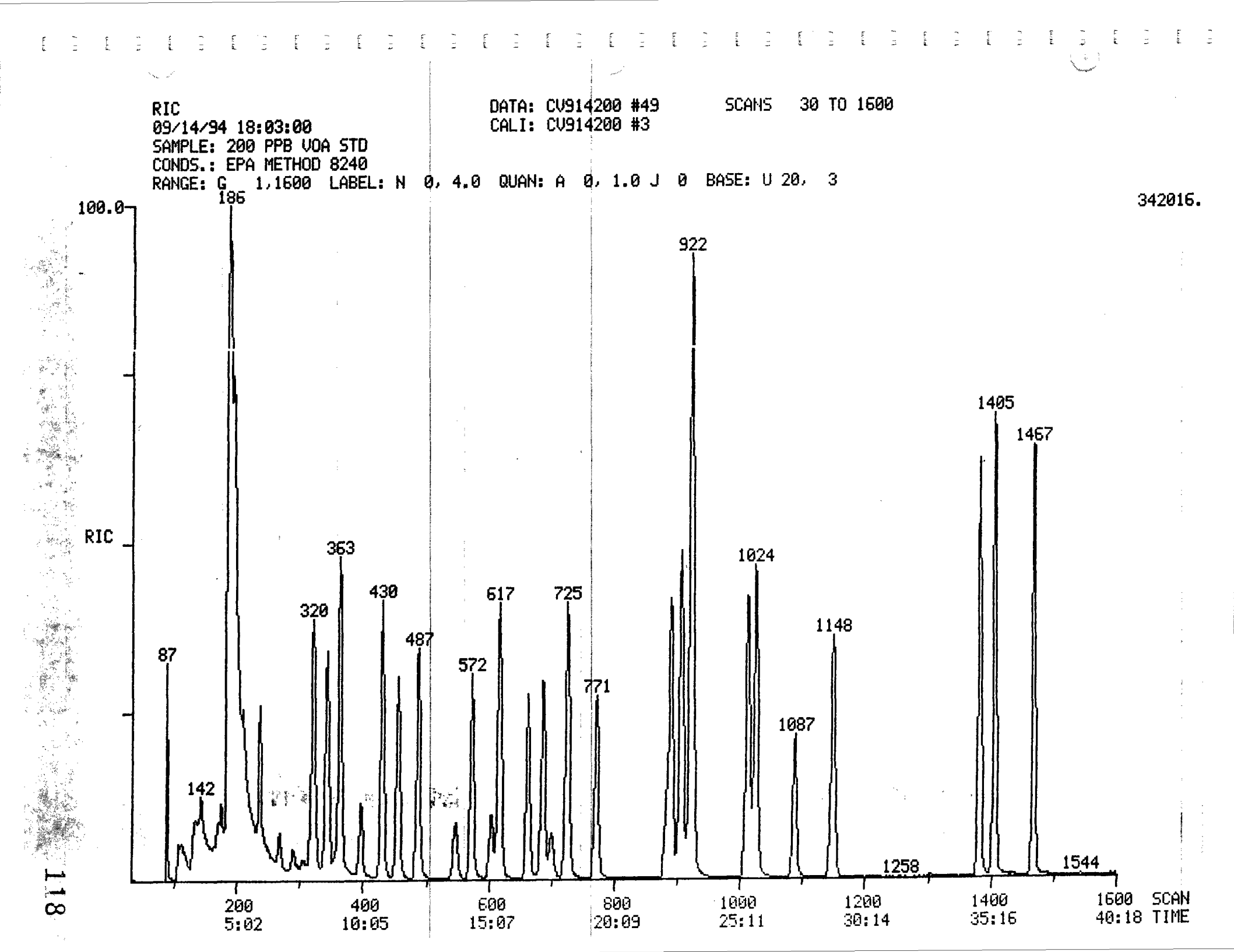
Data: CV914100B.TI
9/14/94 18:53:00
Sample: 100 PPB VO ASTD
Conds.: EPA METHOD 8240
Formula: 5ML Instrument: FINN Weight: 0.000
Submitted by: PTL Analyst: UC Acct. No.: 5-PT

AMOUNT=AREA * REF AMNT/(REF AREA * RESP FACT)
Resp. fac. from Library Entry

No	Name
1	CI01 BROMOCHLOROMETHANE **INT. STD. **
2	CI10 1,4-DIFLUOROBENZENE **INT. STD. **
3	CI20 CHLOROBENZENE-D5 **INT. STD. **
4	CS15 1,2-DICHLOROETHANE-D4 **S. STD. **
5	CS05 TOLUENE-D8 **S. STD. **
6	CS10 4-BROMOFLUOROBENZENE **S. STD. **
7	C010 CHLOROMETHANE **
8	C015 BROMOMETHANE
9	C020 VINYL CHLORIDE *
10	C025 CHLOROETHANE
11	C030 METHYLENE CHLORIDE
12	C251 ACROLIN
13	C035 ACETONE
14	C252 ACRYLONITRILE
15	C040 CARBON DISULFIDE
16	C045 1,1-DICHLOROETHENE *
17	C050 1,1-DICHLOROETHANE **
18	C055 TRANS-1,2-DICHLOROETHENE
19	C000 TRICHLOROFLUOROMETHANE
20	C060 CHLOROFORM *
21	C065 1,2-DICHLOROETHANE
22	C110 2-BUTANONE
23	C115 1,1,1-TRICHLOROETHANE
24	C120 CARBON TETRACHLORIDE
25	C125 VINYL ACETATE
26	C130 BROMO DICHLOROMETHANE
27	C140 1,2-DICHLOROPROPANE *
28	C145 TRANS-1,3 DICHLOROPROPENE
29	C150 TRICHLOROETHENE
30	C155 DIBROMOCHLOROMETHANE
31	C160 1,1,2-TRICHLOROETHANE
32	C165 BENZENE
33	C143 CIS-1,3-DICHLOROPROPENE
34	C175 2-CHLOROETHYL VINYL ETHER
35	C180 BROMOFORM **
36	C220 TETRACHLOROETHENE
37	C210 2-HEXANONE
38	C205 4-METHYL 2-PENTANONE
39	C225 1,1,2,2-TETRACHLOROETHANE **
40	C230 TOLUENE *
41	C235 CHLOROBENZENE **
42	C240 ETHYL BENZENE *
43	C245 STYRENE
44	C250 M+P-XYLENES
45	C253 1,3-DICHLOROBENZENE
46	C254 1,4-DICHLOROBENZENE
47	C255 1,2-DICHLOROBENZENE

No	Name
48	C250 O-XYLENE
49	BENZENE-D6 **S. STD. **

No	m/z	Scan	Time	Ref	RRT	Meth	Area(Hght)	Amount	%Tot
1	49	305	7:41	1	1.000	A BV	8332.	50.000 NG	1.05
2	114	395	9:57	2	1.000	A BB	90012.	50.000 NG	1.05
3	117	879	22:09	3	1.000	A BB	73739.	50.000 NG	1.05
4	65	354	8:55	1	1.161	A BV	13901.	45.807 NG	0.97
5	98	601	15:08	3	0.684	A BB	91027.	50.136 NG	1.06
6	95	1146	28:52	3	1.304	A BB	74302.	52.116 NG	1.10
7	50	107	2:42	1	0.351	A BB	44707.	107.064 NG	2.26
8	94	130	3:16	1	0.426	A BB	50957.	129.774 NG	2.74
9	62	113	2:51	1	0.370	A BB	41066.	94.172 NG	1.99
10	64	133	3:21	1	0.436	A BB	20736.	152.352 NG	3.21
11	49	195	4:55	1	0.639	A BB	53428.	97.468 NG	2.06
12	56	163	4:06	1	0.534	M XX	1978.	148.278 NG	3.13
13	43	166	4:11	1	0.544	M XX	26942.	120.835 NG	2.55
14	53	203	5:07	1	0.666	A BB	10186.	116.630 NG	2.46
15	76	194	4:53	1	0.636	A BB	134740.	107.296 NG	2.26
16	96	171	4:18	1	0.561	A BB	11993.	121.333 NG	2.56
17	63	234	5:54	1	0.767	A BB	12259.	102.063 NG	2.15
18	61	210	5:17	1	0.689	A BB	38757.	108.073 NG	2.28
19	101	142	3:35	1	0.466	A BB	8717.	99.929 NG	2.11
20	83	289	7:17	1	0.948	A BB	15570.	98.942 NG	2.09
21	62	363	9:09	1	1.190	A BB	40391.	104.070 NG	2.19
22	43	266	6:42	2	0.673	A BB	21363.	110.897 NG	2.34
23	97	320	8:04	2	0.810	A BB	130938.	102.154 NG	2.15
24	117	342	8:37	2	0.866	A BB	102726.	104.742 NG	2.21
25	43	235	5:55	2	0.595	A BB	111183.	122.666 NG	2.59
26	83	485	12:13	2	1.228	A BB	125361.	99.993 NG	2.11
27	63	454	11:26	2	1.149	A BB	65276.	101.710 NG	2.15
28	75	570	14:21	2	1.443	A BB	99226.	102.311 NG	2.16
29	95	428	10:47	2	1.084	A BB	78860.	95.245 NG	2.01
30	129	768	19:21	2	1.944	A BB	91532.	102.407 NG	2.16
31	97	683	17:12	2	1.729	A BB	59509.	98.645 NG	2.08
32	78	362	9:07	2	0.916	A BB	187257.	103.006 NG	2.17
33	75	660	16:38	2	1.671	A BB	90090.	103.133 NG	2.18
34	63	541	13:38	2	1.370	M XX	36547.	104.940 NG	2.21
35	173	1085	27:20	2	2.747	A BB	55334.	101.615 NG	2.14
36	166	722	18:11	3	0.821	A BB	72157.	94.697 NG	2.00
37	43	542	13:39	3	0.617	A BB	77799.	106.143 NG	2.24
38	43	696	17:32	3	0.792	A BB	36416.	110.980 NG	2.34
39	83	1145	28:51	3	1.303	A BB	74267.	83.216 NG	1.76
40	91	614	15:28	3	0.699	A BB	207169.	97.186 NG	2.05
41	112	886	22:19	3	1.008	A BB	145047.	99.891 NG	2.11
42	106	903	22:45	3	1.027	A BB	65735.	104.704 NG	2.21
43	104	1022	25:45	3	1.163	A VB	145925.	105.128 NG	2.22
44	106	919	23:09	3	1.046	A BB	176711.	105.374 NG	2.22
45	146	1381	34:47	3	1.571	A BB	126695.	90.753 NG	1.91
46	146	1404	35:22	3	1.597	A BB	126306.	89.172 NG	1.88
47	146	1467	36:57	3	1.669	A BB	113466.	87.482 NG	1.85
48	106	1009	25:25	3	1.148	A BB	87673.	106.802 NG	2.25
49	NOT FOUND								


RIC
09/14/94 18:03:00
SAMPLE: 200 PPB VOA STD
CONDS.: EPA METHOD 8240
RANGE: G 1,1600 LABEL: N 0, 4.0 QUAN: A 0, 1.0 J 0 BASE: U 20, 3
DATA: CV914200 #49 SCANS 30 TO 1600
CALI: CV914200 #3
342016.
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RIC
186
922
1405
1467
363
430
320
617
725
1024
487
1148
87
572
771
1087
142
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200
5:02
400
10:05
600
15:07
800
20:09
1000
25:11
1200
30:14
1400
35:16
1600
40:18
SCAN
TIME
118

Quantitation Report File: CV914200

Data: CV914200.TI
)/14/94 18:03:00
Sample: 200 PPB VOA STD
Conds.: EPA METHOD 8240
Formula: 5ML Instrument: FINN Weight: 0.000
Submitted by: PTL Analyst: UC Acct. No.: 5-PT

AMOUNT=AREA * REF AMNT/(REF AREA * RESP FACT)
Resp. fac. from Library Entry

No	Name
1	CI01 BROMOCHLOROMETHANE **INT. STD. **
2	CI10 1,4-DIFLUOROBENZENE **INT. STD. **
3	CI20 CHLOROBENZENE-D5 **INT. STD. **
4	CS15 1,2-DICHLOROETHANE-D4 **S. STD. **
5	CS05 TOLUENE-D8 **S. STD. **
6	CS10 4-BROMOFLUOROBENZENE **S. STD. **
7	C010 CHLOROMETHANE **
8	C015 BROMOMETHANE
9	C020 VINYL CHLORIDE *
10	C025 CHLOROETHANE
11	C030 METHYLENE CHLORIDE
12	C251 ACROLIN
13	C035 ACETONE
14	C252 ACRYLONITRILE
15	C040 CARBON DISULFIDE
16	C045 1,1-DICHLOROETHENE *
17	C050 1,1-DICHLOROETHANE **
18	C055 TRANS-1,2-DICHLOROETHENE
19	C000 TRICHLOROFLUOROMETHANE
20	C060 CHLOROFORM *
21	C065 1,2-DICHLOROETHANE
22	C110 2-BUTANONE
23	C115 1,1,1-TRICHLOROETHANE
24	C120 CARBON TETRACHLORIDE
25	C125 VINYL ACETATE
26	C130 BROMO DICHLOROMETHANE
27	C140 1,2-DICHLOROPROPANE *
28	C145 TRANS-1,3 DICHLOROPROPENE
29	C150 TRICHLOROETHENE
30	C155 DIBROMOCHLOROMETHANE
31	C160 1,1,2-TRICHLOROETHANE
32	C165 BENZENE
33	C143 CIS-1,3-DICHLOROPROPENE
34	C175 2-CHLOROETHYL VINYL ETHER
35	C180 BROMOFORM **
36	C220 TETRACHLOROETHENE
37	C210 2-HEXANONE
38	C205 4-METHYL 2-PENTANONE
39	C225 1,1,2,2-TETRACHLOROETHANE **
40	C230 TOLUENE *
41	C235 CHLOROBENZENE **
42	C240 ETHYL BENZENE *
43	C245 STYRENE
44	C250 M+P-XYLENES
45	C253 1,3-DICHLOROBENZENE
46	C254 1,4-DICHLOROBENZENE
47	C255 1,2-DICHLOROBENZENE

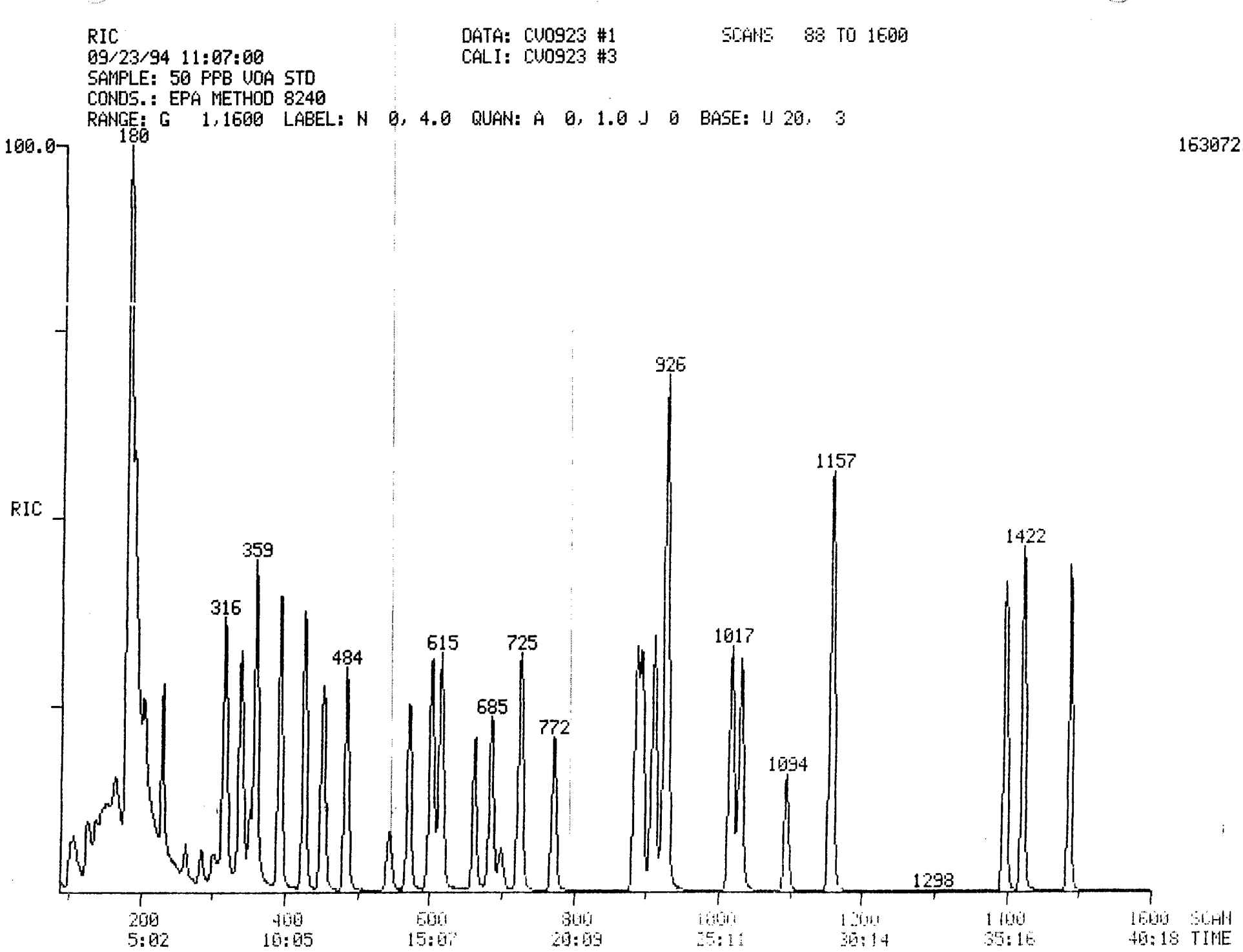
No	Name
48	C250 O-XYLENE
49	BENZENE-D6 **S. STD. **

No	m/z	Scan	Time	Ref	RRT	Meth	Area(Hght)	Amount	%Tot
1	49	306	7:42	1	1.000	A BV	7806.	50.000 NG	0.52
2	114	396	9:59	2	1.000	A BB	88807.	50.000 NG	0.52
3	117	882	22:13	3	1.000	A BB	75395.	50.000 NG	0.52
4	65	355	8:57	1	1.160	A BV	13515.	47.539 NG	0.50
5	98	603	15:11	3	0.684	A BB	92800.	49.991 NG	0.52
6	95	1148	28:55	3	1.302	A BB	102798.	70.521 NG	0.74
7	50	107	2:42	1	0.350	A BB	69405.	177.418 NG	1.86
8	94	131	3:18	1	0.428	A BB	84863.	230.699 NG	2.42
9	62	112	2:49	1	0.366	A BB	52850.	129.369 NG	1.36
10	64	135	3:24	1	0.441	M XX	36061.	282.812 NG	2.97
11	49	195	4:55	1	0.637	A BB	121356.	236.319 NG	2.48
12	56	164	4:08	1	0.536	A BV	8317.	665.676 NG	6.98
13	43	175	4:24	1	0.572	M XX	41602.	199.167 NG	2.09
14	53	205	5:10	1	0.670	A BB	17717.	216.540 NG	2.27
15	76	194	4:53	1	0.634	A BB	293321.	249.329 NG	2.62
16	96	174	4:23	1	0.569	A BB	25442.	274.760 NG	2.88
17	63	235	5:55	1	0.768	A BB	26555.	235.996 NG	2.48
18	61	209	5:16	1	0.683	A BB	78604.	233.963 NG	2.45
19	101	142	3:35	1	0.464	A BB	22594.	276.477 NG	2.90
20	83	289	7:17	1	0.944	A BB	36333.	246.447 NG	2.59
21	62	364	9:10	1	1.190	A BB	82722.	227.512 NG	2.39
22	43	267	6:44	2	0.674	A BB	51123.	268.984 NG	2.82
23	97	320	8:04	2	0.808	A BB	269582.	213.173 NG	2.24
24	117	342	8:37	2	0.864	A BB	214553.	221.732 NG	2.33
25	43	235	5:55	2	0.593	A BB	207998.	232.594 NG	2.44
26	83	487	12:16	2	1.230	A BB	257963.	208.554 NG	2.19
27	63	455	11:28	2	1.149	A BB	130599.	206.255 NG	2.16
28	75	572	14:25	2	1.444	A BB	201586.	210.673 NG	2.21
29	95	430	10:50	2	1.086	A BB	160149.	196.048 NG	2.06
30	129	771	19:25	2	1.947	A BB	185177.	209.989 NG	2.20
31	97	685	17:15	2	1.730	A BB	121282.	203.771 NG	2.14
32	78	362	9:07	2	0.914	A BB	359869.	200.641 NG	2.11
33	75	662	16:41	2	1.672	A BB	176297.	204.561 NG	2.15
34	NOT FOUND								
35	173	1087	27:23	2	2.745	A BB	126297.	235.078 NG	2.47
36	166	725	18:16	3	0.822	A BB	150600.	193.304 NG	2.03
37	43	546	13:45	3	0.619	A BB	87537.	116.807 NG	1.23
38	43	699	17:36	3	0.793	A BB	68826.	205.148 NG	2.15
39	83	1147	28:54	3	1.300	A BB	187103.	205.045 NG	2.15
40	91	616	15:31	3	0.698	A BB	424538.	194.784 NG	2.04
41	112	889	22:24	3	1.008	A BB	294155.	198.131 NG	2.08
42	106	905	22:48	3	1.026	A BB	136028.	211.911 NG	2.22
43	104	1024	25:48	3	1.161	A VB	311960.	219.809 NG	2.31
44	106	922	23:14	3	1.045	A BB	339687.	198.111 NG	2.08
45	146	1382	34:49	3	1.567	A BB	272771.	191.099 NG	2.01
46	146	1405	35:24	3	1.593	A BB	275728.	190.392 NG	2.00
47	146	1467	36:57	3	1.663	A BB	255873.	192.945 NG	2.02
48	106	1011	25:28	3	1.146	A BB	168185.	200.384 NG	2.10
49	NOT FOUND								

DATAFILE: CV0923 Calibration Check

Instrument Identifier: FINN
Calibration Date: 09/14/94
Standard File: CV0923
Date: 09/23/94 Time: 11:07:00
25% D
MIN RF FOR SPCC (**) = 0.300

	Compound	Mean RF(I)	RF(0)	% D
CS15	1,2-DICHLOROETHANE-D4 *	2.017	2.528	0.000
CS05	TOLUENE-D8 **S. STD. **	1.232	1.339	0.000
CS10	4-BROMOFLUOROBENZENE **	1.020	1.035	0.000
C010	CHLOROMETHANE **	2.312	2.971	28.489
C015	BROMOMETHANE	2.485	2.560	3.036
C020	VINYL CHLORIDE *	2.413	2.186	9.414
C025	CHLOROETHANE	0.920	0.714	22.364
C030	METHYLENE CHLORIDE	3.720	3.623	2.599
C251	ACROLIN	0.190	0.547	187.463
C035	ACETONE	1.417	1.737	22.628
C252	ACRYLONITRILE	0.513	0.643	25.311
C040	CARBON DISULFIDE	8.003	9.790	22.337
C045	1,1-DICHLOROETHENE *	0.678	0.733	8.243
C050	1,1-DICHLOROETHANE **	0.807	1.192	47.796
C055	TRANS-1,2-DICHLOROETHEN	2.272	2.987	31.471
C000	TRICHLOROFLUOROMETHA	0.601	0.586	2.414
C060	CHLOROFORM *	1.124	1.329	18.213
C065	1,2-DICHLOROETHANE	2.724	3.305	21.309
C110	2-BUTANONE	0.133	0.148	11.402
C115	1,1,1-TRICHLOROETHANE	0.711	0.748	5.236
C120	CARBON TETRACHLORIDE	0.561	0.601	7.108
C125	VINYL ACETATE	0.466	0.691	48.160
C130	BROMO DICHLOROMETHANE	0.668	0.689	3.235
C140	1,2-DICHLOROPROPANE	0.352	0.356	1.218
C145	TRANS-1,3 DICHLOROPROPE	0.510	0.489	4.124
C150	TRICHLOROETHENE	0.440	0.435	1.303
C155	DIBROMOCHLOROMETHANE	0.482	0.429	11.015
C160	1,1,2-TRICHLOROETHANE	0.329	0.288	12.256
C165	BENZENE	0.994	0.959	3.583
C143	CIS-1,3-DICHLOROPROPENE	0.458	0.392	14.322
C175	2-CHLOROETHYL VINYL ETH	0.144	0.015	89.347
C180	BROMOFORM **	0.306	0.285	6.792
C220	TETRACHLOROETHENE	0.501	0.523	4.394
C210	2-HEXANONE	0.449	0.390	13.127
C205	4-METHYL 2-PENTANONE	0.249	0.278	11.505
C225	1,1,2,2-TETRACHLOROETHA	0.596	0.685	14.898
C230	TOLUENE *	1.392	1.466	5.331
C235	CHLOROBENZENE **	0.968	0.937	3.226
C240	ETHYL BENZENE *	0.424	0.408	3.785
C245	STYRENE	0.925	0.944	2.036
C250	M+P-XYLENES	1.102	1.167	5.884
C253	1,3-DICHLOROBENZENE	0.899	0.904	0.586
C254	1,4-DICHLOROBENZENE	0.909	0.906	0.369
C255	1,2-DICHLOROBENZENE	0.850	0.816	4.068
C250	O-XYLENE	0.559	0.587	4.990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09/23/94 11:07:00
SAMPLE: 50 PPB VOA STD
CONDS.: EPA METHOD 8240
RANGE: G 1,1600 LABEL: N 0, 4.0 QUAN: A 0, 1.0 J 0 BASE: U 20, 3
DATA: CV0923 #1
CALI: CV0923 #3
SCANS 88 TO 1600
100.0
163072.
RIC
180
316
359
484
615
685
725
772
926
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1094
1157
1298
1422
200
5:02
400
10:05
600
15:07
800
20:09
1000
25:11
1200
30:14
1400
35:16
1600 SCAN
40:18 TIME

Quantitation Report File: CV0923

Data: CV0923.TI
9/23/94 11:07:00
Sample: 50 PPB VOA STD
Conds.: EPA METHOD 8240
Formula: 5ML Instrument: FINN Weight: 0.000
Submitted by: PTL Analyst: UC Acct. No.:

AMOUNT=AREA * REF AMNT/(REF AREA * RESP FACT)
Resp. fac. from Library Entry

No	Name
1	CI01 BROMOCHLOROMETHANE **INT. STD. **
2	CI10 1,4-DIFLUOROBENZENE **INT. STD. **
3	CI20 CHLOROBENZENE-D5 **INT. STD. **
4	CS15 1,2-DICHLOROETHANE-D4 **S. STD. **
5	CS05 TOLUENE-D8 **S. STD. **
6	CS10 4-BROMOFLUOROBENZENE **S. STD. **
7	C010 CHLOROMETHANE **
8	C015 BROMOMETHANE
9	C020 VINYL CHLORIDE *
10	C025 CHLOROETHANE
11	C030 METHYLENE CHLORIDE
12	C251 ACROLIN
13	C035 ACETONE
14	C252 ACRYLONITRILE
15	C040 CARBON DISULFIDE
16	C045 1,1-DICHLOROETHENE *
17	C050 1,1-DICHLOROETHANE **
18	C055 TRANS-1,2-DICHLOROETHENE
19	C000 TRICHLOROFLUOROMETHANE
20	C060 CHLOROFORM *
21	C065 1,2-DICHLOROETHANE
22	C110 2-BUTANONE
23	C115 1,1,1-TRICHLOROETHANE
24	C120 CARBON TETRACHLORIDE
25	C125 VINYL ACETATE
26	C130 BROMO DICHLOROMETHANE
27	C140 1,2-DICHLOROPROPANE *
28	C145 TRANS-1,3 DICHLOROPROPENE
29	C150 TRICHLOROETHENE
30	C155 DIBROMOCHLOROMETHANE
31	C160 1,1,2-TRICHLOROETHANE
32	C165 BENZENE
33	C143 CIS-1,3-DICHLOROPROPENE
34	C175 2-CHLOROETHYL VINYL ETHER
35	C180 BROMOFORM **
36	C220 TETRACHLOROETHENE
37	C210 2-HEXANONE
38	C205 4-METHYL 2-PENTANONE
39	C225 1,1,2,2-TETRACHLOROETHANE **
40	C230 TOLUENE *
41	C235 CHLOROBENZENE **
42	C240 ETHYL BENZENE *
43	C245 STYRENE
44	C250 M+P-XYLENES
45	C253 1,3-DICHLOROBENZENE
46	C254 1,4-DICHLOROBENZENE
47	C255 1,2-DICHLOROBENZENE

(W923

No	Name
48	C250 O-XYLENE
49	BENZENE-D6 **S. STD. **

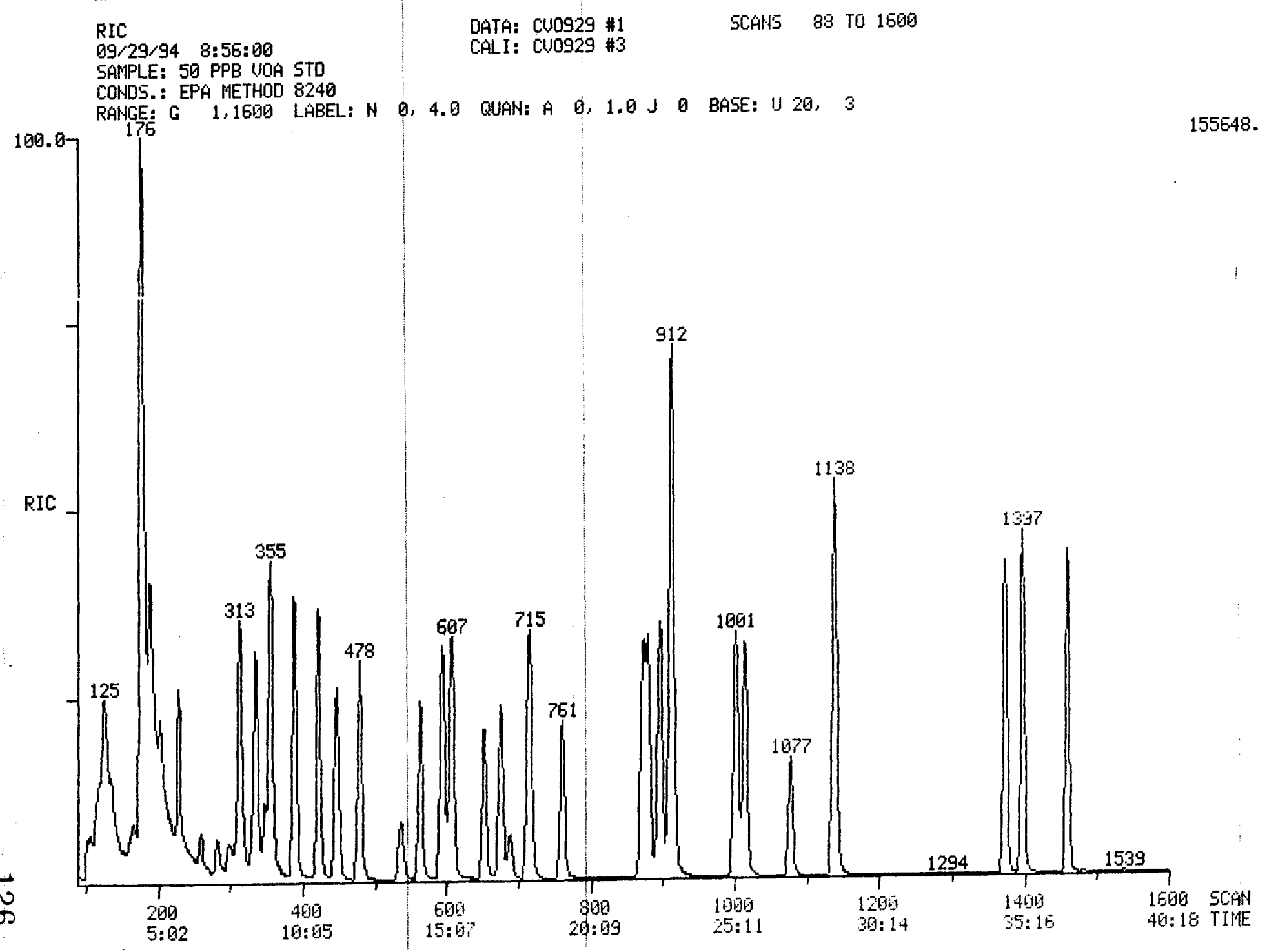
No	m/z	Scan	Time	Ref	RRT	Meth	Area(Hght)	Amount	%Tot
1	49	301	7:35	1	1.000	A BV	12697.	50.000 NG	1.89
2	114	392	9:52	2	1.000	A BB	158172.	50.000 NG	1.89
3	117	885	22:18	3	1.000	A BB	106406.	50.000 NG	1.89
4	65	350	8:49	1	1.163	A BB	32098.	75.118 NG	2.85
5	98	601	15:08	3	0.679	A BB	142422.	50.830 NG	1.93
6	95	1157	29:09	3	1.307	A BB	110045.	50.254 NG	1.90
7	50	102	2:34	1	0.339	A BB	37716.	81.895 NG	3.10
8	94	125	3:09	1	0.415	M XX	32502.	33.746 NG	1.28
9	62	107	2:42	1	0.355	A BB	27754.	71.606 NG	2.71
10	64	128	3:13	1	0.425	A BB	9065.	29.633 NG	1.12
11	49	190	4:47	1	0.631	A BB	45995.	62.920 NG	2.38
12	56	159	4:00	1	0.528	M XX	6936.	100.602 NG	3.81
13	43	162	4:05	1	0.538	M XX	29646.	99.713 NG	3.78
14	53	199	5:01	1	0.661	A BB	8155.	71.113 NG	2.69
15	76	189	4:46	1	0.628	A BB	124308.	66.442 NG	2.52
16	96	165	4:09	1	0.548	A BB	9307.	61.837 NG	2.34
17	63	229	5:46	1	0.761	A BB	15133.	76.189 NG	2.89
18	61	205	5:10	1	0.681	A BB	37915.	90.428 NG	3.43
19	101	136	3:26	1	0.452	A BB	7439.	18.782 NG	0.71
20	83	284	7:09	1	0.944	A BB	16862.	62.470 NG	2.37
21	62	360	9:04	1	1.196	A BB	41953.	76.491 NG	2.90
22	43	262	6:36	2	0.668	A BB	23352.	55.556 NG	2.11
23	97	316	7:58	2	0.806	A BB	118217.	51.396 NG	1.95
24	117	338	8:31	2	0.862	A BB	95005.	51.509 NG	1.95
25	43	230	5:48	2	0.587	A BB	109147.	59.098 NG	2.24
26	83	484	12:12	2	1.235	A BB	108921.	49.332 NG	1.87
27	63	451	11:22	2	1.151	A BB	56281.	51.507 NG	1.95
28	75	571	14:23	2	1.457	A BB	77210.	49.924 NG	1.89
29	95	426	10:44	2	1.087	A BB	68666.	49.133 NG	1.86
30	129	772	19:27	2	1.969	A BB	67793.	41.771 NG	1.58
31	97	685	17:15	2	1.747	A BB	45517.	42.502 NG	1.61
32	78	358	9:01	2	0.913	A BB	151586.	49.225 NG	1.87
33	75	662	16:41	2	1.689	A BB	61968.	47.574 NG	1.80
34	63	541	13:38	2	1.380	M XX	2353.	41.846 NG	1.59
35	173	1094	27:33	2	2.791	A BB	45060.	38.430 NG	1.46
36	166	725	18:16	3	0.819	A BB	55565.	40.081 NG	1.52
37	43	543	13:41	3	0.614	A BB	41430.	61.146 NG	2.32
38	43	699	17:36	3	0.790	A BB	29494.	59.870 NG	2.27
39	83	1156	29:07	3	1.306	A BB	72803.	48.981 NG	1.86
40	91	615	15:30	3	0.695	A BB	155960.	49.191 NG	1.86
41	112	893	22:30	3	1.009	A BB	99663.	48.438 NG	1.84
42	106	910	22:55	3	1.028	A BB	43357.	50.360 NG	1.91
43	104	1031	25:58	3	1.165	A VB	100359.	51.435 NG	1.95
44	106	926	23:20	3	1.046	A BB	124095.	49.812 NG	1.89
45	146	1397	35:11	3	1.579	A BB	96150.	40.664 NG	1.54
46	146	1422	35:49	3	1.607	A BB	96364.	39.443 NG	1.49
47	146	1487	37:27	3	1.680	A BB	86757.	39.822 NG	1.51
48	106	1017	25:37	3	1.149	A BB	62380.	50.776 NG	1.92
49	NOT FOUND								

DATAFILE: CV0929

Calibration Check

Instrument Identifier: FINN
Calibration Date: 09/14/94
Standard File: CV0929
Date: 09/29/94 Time: 08:56:00
25% D
MIN RF FOR SPCC (**) = 0.300

	Compound	Mean RF(I)	RF(0)	% D
CS15	1,2-DICHLOROETHANE-D4 *	2.017	1.692	0.000
CS05	TOLUENE-D8 **S. STD. **	1.232	1.340	0.000
CS10	4-BROMOFLUOROBENZENE **	1.020	0.969	0.000
C010	CHLOROMETHANE **	2.312	1.942	16.015
C015	BROMOMETHANE	2.485	3.271	31.658
C020	VINYL CHLORIDE *	2.413	2.388	1.073
C025	CHLOROETHANE	0.920	1.145	24.396
C030	METHYLENE CHLORIDE	3.720	2.442	34.336
C251	ACROLIN	0.190	0.096	49.341
C035	ACETONE	1.417	0.990	30.147
C252	ACRYLONITRILE	0.513	0.448	12.592
C040	CARBON DISULFIDE	8.003	7.054	11.859
C045	1,1-DICHLOROETHENE *	0.678	0.549	18.916
C050	1,1-DICHLOROETHANE **	0.807	1.092	35.310
C055	TRANS-1,2-DICHLOROETHEN	2.272	1.779	21.668
C000	TRICHLOROFLUOROMETHA	0.601	0.546	9.158
C060	CHLOROFORM *	1.124	1.079	3.999
C065	1,2-DICHLOROETHANE	2.724	2.269	16.724
110	2-BUTANONE	0.133	0.153	15.353
C115	1,1,1-TRICHLOROETHANE	0.711	0.735	3.486
C120	CARBON TETRACHLORIDE	0.561	0.594	5.892
C125	VINYL ACETATE	0.466	0.696	49.265
C130	BROMO DICHLOROMETHANE	0.668	0.703	5.373
C140	1,2-DICHLOROPROPANE	0.352	0.364	3.367
C145	TRANS-1,3 DICHLOROPROPE	0.510	0.512	0.516
C150	TRICHLOROETHENE	0.440	0.428	2.797
C155	DIBROMOCHLOROMETHANE	0.482	0.466	3.415
C160	1,1,2-TRICHLOROETHANE	0.329	0.314	4.311
C165	BENZENE	0.994	0.982	1.259
C143	CIS-1,3-DICHLOROPROPENE	0.458	0.433	5.528
C175	2-CHLOROETHYL VINYL ETH	0.144	0.017	87.994
C180	BROMOFORM **	0.306	0.329	7.321
C220	TETRACHLOROETHENE	0.501	0.593	18.461
C210	2-HEXANONE	0.449	0.383	14.689
C205	4-METHYL 2-PENTANONE	0.249	0.277	11.134
C225	1,1,2,2-TETRACHLOROETHA	0.596	0.676	13.414
C230	TOLUENE *	1.392	1.549	11.273
C235	CHLOROBENZENE **	0.968	0.962	0.670
C240	ETHYL BENZENE *	0.424	0.423	0.218
C245	STYRENE	0.925	0.984	6.402
C250	M+P-XYLENES	1.102	1.200	8.908
C253	1,3-DICHLOROBENZENE	0.899	0.934	3.967
C254	1,4-DICHLOROBENZENE	0.909	0.932	2.522
C255	1,2-DICHLOROBENZENE	0.850	0.820	3.624
C250	O-XYLENE	0.559	0.585	4.663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09/29/94 8:56:00
SAMPLE: 50 PPB VOA STD
CONDS.: EPA METHOD 8240
RANGE: G 1,1600 LABEL: N 0, 4.0 QUAN: A 0, 1.0 J 0 BASE: U 20, 3
DATA: CV0929 #1
CALI: CV0929 #3
SCANS 88 TO 1600
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100.0
RIC
176
125
313
355
478
607
715
761
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1001
1077
1138
1294
1397
1539
200
5:02
400
10:05
600
15:07
800
20:09
1000
25:11
1200
30:14
1400
35:16
1600 SCAN
40:18 TIME

Quantitation Report File: CV0929

Data: CV0929.TI
9/29/94 8:56:00
Sample: 50 PPB VOA STD
Conds.: EPA METHOD 8240
Formula: 5ML Instrument: FINN Weight: 0.000
Submitted by: PTL Analyst: UC Acct. No.:

AMOUNT=AREA * REF AMNT/(REF AREA * RESP FACT)
Resp. fac. from Library Entry

No	Name
1	CI01 BROMOCHLOROMETHANE **INT. STD. **
2	CI10 1,4-DIFLUOROBENZENE **INT. STD. **
3	CI20 CHLOROBENZENE-D5 **INT. STD. **
4	CS15 1,2-DICHLOROETHANE-D4 **S. STD. **
5	CS05 TOLUENE-D8 **S. STD. **
6	CS10 4-BROMOFLUOROBENZENE **S. STD. **
7	C010 CHLOROMETHANE **
8	C015 BROMOMETHANE
9	C020 VINYL CHLORIDE *
10	C025 CHLOROETHANE
11	C030 METHYLENE CHLORIDE
12	C251 ACROLIN
13	C035 ACETONE
14	C252 ACRYLONITRILE
15	C040 CARBON DISULFIDE
16	C045 1,1-DICHLOROETHENE *
17	C050 1,1-DICHLOROETHANE **
18	C055 TRANS-1,2-DICHLOROETHENE
19	C000 TRICHLOROFLUOROMETHANE
20	C060 CHLOROFORM *
21	C065 1,2-DICHLOROETHANE
22	C110 2-BUTANONE
23	C115 1,1,1-TRICHLOROETHANE
24	C120 CARBON TETRACHLORIDE
25	C125 VINYL ACETATE
26	C130 BROMO DICHLOROMETHANE
27	C140 1,2-DICHLOROPROPANE *
28	C145 TRANS-1,3 DICHLOROPROPENE
29	C150 TRICHLOROETHENE
30	C155 DIBROMOCHLOROMETHANE
31	C160 1,1,2-TRICHLOROETHANE
32	C165 BENZENE
33	C143 CIS-1,3-DICHLOROPROPENE
34	C175 2-CHLOROETHYL VINYL ETHER
35	C180 BROMOFORM **
36	C220 TETRACHLOROETHENE
37	C210 2-HEXANONE
38	C205 4-METHYL 2-PENTANONE
39	C225 1,1,2,2-TETRACHLOROETHANE **
40	C230 TOLUENE *
41	C235 CHLOROBENZENE **
42	C240 ETHYL BENZENE *
43	C245 STYRENE
44	C250 M+P-XYLENES
45	C253 1,3-DICHLOROBENZENE
46	C254 1,4-DICHLOROBENZENE
47	C255 1,2-DICHLOROBENZENE

Cno 929

No	Name
48	C250 O-XYLENE
49	BENZENE-D6 **S. STD. **

No	m/z	Scan	Time	Ref	RRT	Meth	Area(Hght)	Amount	%Tot
1	49	298	7:30	1	1.000	M XX	17120.	50.000 NG	2.08
2	114	388	9:46	2	1.000	A BB	142012.	50.000 NG	2.08
3	117	872	21:58	3	1.000	A BB	99603.	50.000 NG	2.08
4	65	347	8:44	1	1.164	A BB	28952.	50.000 NG	2.08
5	98	594	14:58	3	0.681	A BB	133468.	50.000 NG	2.08
6	95	1139	28:41	3	1.306	A BB	96440.	50.000 NG	2.08
7	50	102	2:34	1	0.342	A BB	33236.	50.000 NG	2.08
8	94	125	3:09	1	0.419	M XX	55998.	50.000 NG	2.08
9	62	107	2:42	1	0.359	M XX	40867.	50.000 NG	2.08
10	64	128	3:13	1	0.430	M XX	19590.	50.000 NG	2.08
11	49	189	4:46	1	0.634	A BB	41806.	50.000 NG	2.08
12	56	158	3:59	1	0.530	A BB	1641.	50.000 NG	2.08
13	43	160	4:02	1	0.537	M XX	16933.	50.000 NG	2.08
14	53	198	4:59	1	0.664	A BB	7667.	50.000 NG	2.08
15	76	189	4:46	1	0.634	A BB	120753.	50.000 NG	2.08
16	96	165	4:09	1	0.554	A BB	9398.	50.000 NG	2.08
17	63	228	5:45	1	0.765	M XX	18680.	50.000 NG	2.08
18	61	203	5:07	1	0.681	A BB	30455.	50.000 NG	2.08
19	101	136	3:26	1	0.456	A BB	9336.	50.000 NG	2.08
20	83	282	7:06	1	0.946	A BB	18462.	50.000 NG	2.08
21	62	357	9:00	1	1.198	A BB	38828.	50.000 NG	2.08
22	43	259	6:31	2	0.668	A BB	21712.	50.000 NG	2.08
23	97	313	7:53	2	0.807	A BB	104374.	50.000 NG	2.08
24	117	335	8:26	2	0.863	A BB	84330.	50.000 NG	2.08
25	43	229	5:46	2	0.590	M XX	98727.	50.000 NG	2.08
26	83	478	12:02	2	1.232	A BB	99820.	50.000 NG	2.08
27	63	446	11:14	2	1.149	A BB	51605.	50.000 NG	2.08
28	75	563	14:11	2	1.451	A BB	72680.	50.000 NG	2.08
29	95	421	10:36	2	1.085	A BB	60716.	50.000 NG	2.08
30	129	761	19:10	2	1.961	A BB	66071.	50.000 NG	2.08
31	97	676	17:02	2	1.742	A BB	44574.	50.000 NG	2.08
32	78	355	8:57	2	0.915	A BB	139382.	50.000 NG	2.08
33	75	652	16:25	2	1.680	A BB	61355.	50.000 NG	2.08
34	63	533	13:26	2	1.374	M XX	2390.	50.000 NG	2.08
35	173	1077	27:08	2	2.776	A BB	46593.	50.000 NG	2.08
36	166	715	18:01	3	0.820	A BB	59028.	50.000 NG	2.08
37	43	536	13:30	3	0.615	A BB	38083.	50.000 NG	2.08
38	43	689	17:21	3	0.790	A BB	27516.	50.000 NG	2.08
39	83	1138	28:40	3	1.305	M XX	67268.	50.000 NG	2.08
40	91	607	15:17	3	0.696	A BB	154227.	50.000 NG	2.08
41	112	879	22:09	3	1.008	A BB	95756.	50.000 NG	2.08
42	106	896	22:34	3	1.028	A BB	42092.	50.000 NG	2.08
43	104	1014	25:33	3	1.163	A VB	97966.	50.000 NG	2.08
44	106	912	22:58	3	1.046	A BB	119481.	50.000 NG	2.08
45	146	1373	34:35	3	1.575	A BB	93029.	50.000 NG	2.08
46	146	1397	35:11	3	1.602	A BB	92822.	50.000 NG	2.08
47	146	1460	36:47	3	1.674	A BB	81587.	50.000 NG	2.08
48	106	1001	25:13	3	1.148	A BB	58210.	50.000 NG	2.08
49	NOT FOUND								

DATAFILE: CV0104 Calibration Check

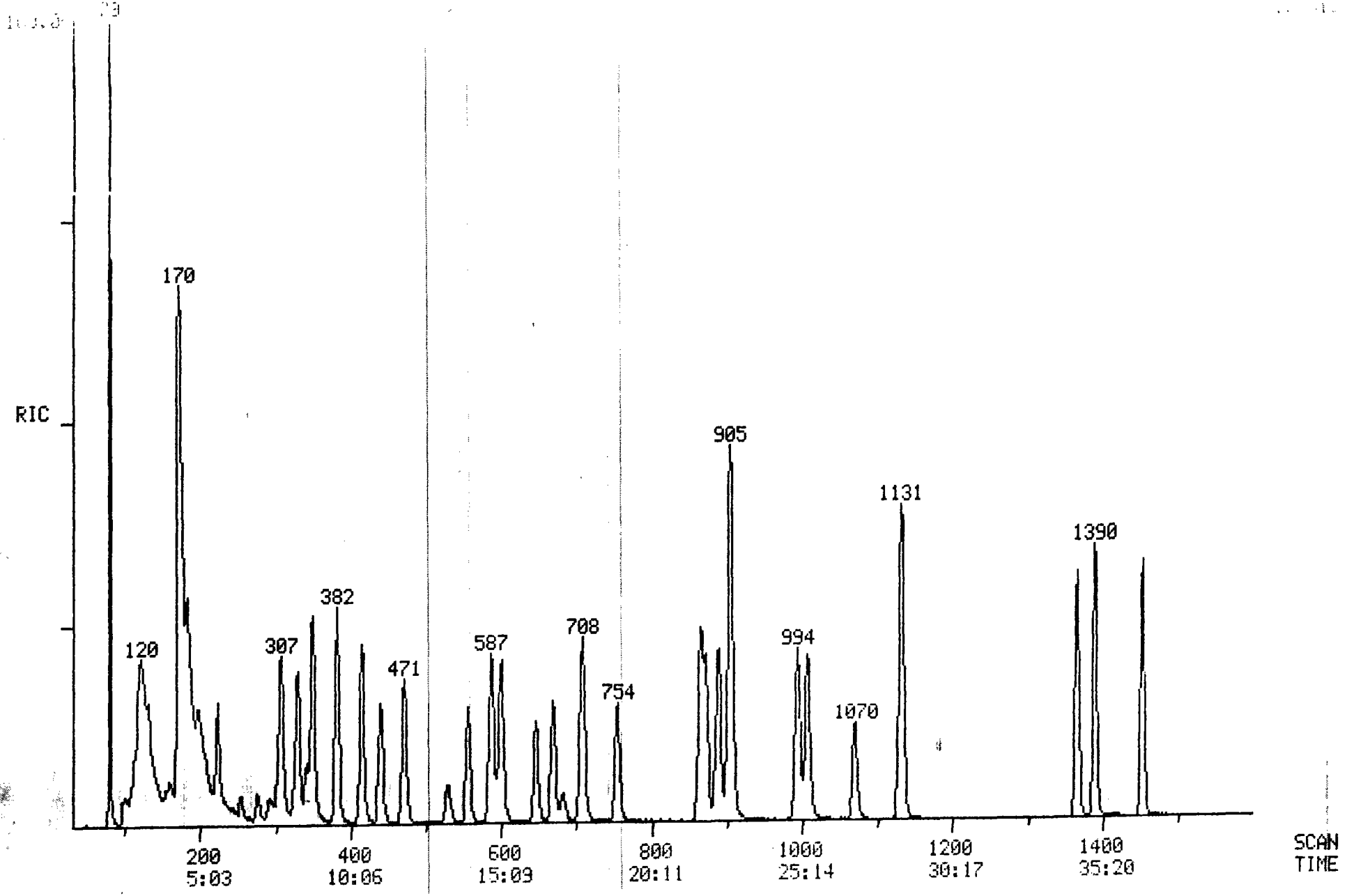
Instrument Identifier: FINN
Calibration Date: 09/14/94
Standard File: CV0104
Date: 10/04/94 Time: 12:34:00
25% D
MIN RF FOR SPCC (**) = 0.300

	Compound	Mean RF(I)	RF(0)	% D
CS15	1,2-DICHLOROETHANE-D4 *	2.017	1.424	0.000
CS05	TOLUENE-D8 **S. STD. **	1.232	1.226	0.000
CS10	4-BROMOFLUOROBENZENE **	1.020	0.939	0.000
C010	CHLOROMETHANE **	2.312	1.160	49.825
C015	BROMOMETHANE	2.485	3.164	27.344
C020	VINYL CHLORIDE *	2.413	1.987	17.663
C025	CHLOROETHANE	0.920	1.598	73.625
C030	METHYLENE CHLORIDE	3.720	1.884	49.361
C251	ACROLIN	0.190	0.310	63.029
C038	ACETONE	1.417	0.930	34.357
C252	ACRYLONITRILE	0.513	0.415	19.048
C040	CARBON DISULFIDE	8.003	5.123	35.983
C045	1,1-DICHLOROETHENE *	0.678	0.456	32.766
C050	1,1-DICHLOROETHANE **	0.807	0.860	6.593
C055	TRANS-1,2-DICHLOROETHEN	2.272	1.593	29.892
C000	TRICHLOROFLUOROMETHA	0.601	0.608	1.222
C060	CHLOROFORM *	1.124	0.849	24.467
C065	1,2-DICHLOROETHANE	2.724	1.762	35.334
C110	2-BUTANONE	0.133	0.123	7.555
C115	1,1,1-TRICHLOROETHANE	0.711	0.615	13.430
C120	CARBON TETRACHLORIDE	0.561	0.512	8.801
C125	VINYL ACETATE	0.466	0.527	12.984
C130	BROMO DICHLOROMETHANE	0.668	0.605	9.353
C140	1,2-DICHLOROPROPANE	0.352	0.289	17.838
C145	TRANS-1,3 DICHLOROPROPE	0.510	0.421	17.419
C150	TRICHLOROETHENE	0.440	0.375	14.780
C155	DIBROMOCHLOROMETHANE	0.482	0.451	6.551
C160	1,1,2-TRICHLOROETHANE	0.329	0.282	14.060
C165	BENZENE	0.994	0.804	19.191
C143	CIS-1,3-DICHLOROPROPENE	0.458	0.354	22.676
C175	2-CHLOROETHYL VINYL ETH	0.144	0.020	85.807
C180	BROMOFORM **	0.306	0.335	9.358
C220	TETRACHLOROETHENE	0.501	0.568	13.511
C210	2-HEXANONE	0.449	0.284	36.811
C205	4-METHYL 2-PENTANONE	0.249	0.212	14.731
C225	1,1,2,2-TETRACHLOROETHA	0.596	0.627	5.186
C230	TOLUENE *	1.392	1.281	7.960
C235	CHLOROBENZENE **	0.968	0.824	14.950
C240	ETHYL BENZENE *	0.424	0.355	16.274
C245	STYRENE	0.925	0.840	9.181
C250	M+P-XYLENES	1.102	1.060	3.819
C253	1,3-DICHLOROBENZENE	0.899	0.943	4.960
C254	1,4-DICHLOROBENZENE	0.909	0.957	5.176
C255	1,2-DICHLOROBENZENE	0.850	0.853	0.313
C250	O-XYLENE	0.559	0.541	3.242

RIC
10/04/94 12:34:00
SAMPLE: 50 PPB VOA STD
CONDS.: EPA METHOD 8240

DATA: CV0104 #49
CALI: CV0104 #3

SCANS 30 TO 1597



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Quantitat: eport File: CV0104

Data: CV0
0/04/94 00
Sample: S VOA STD
Conds.: EP THOD 8240
Formula: Instrument: FINN Weight: 0.000
Submitted TL Analyst: UC Acct. No.: -

AMOUNT=ARE EF AMNT/(REF AREA * RESP FACT)
Resp. fac n Library Entry

No	Name
1	CI01 DICHLOROMETHANE **INT. STD. **
2	CI10 -DIFLUOROBENZENE **INT. STD. **
3	CI20 OBENZENE-D5 **INT. STD. **
4	CS15 ICHLOROETHANE-D4 **S. STD. **
5	CS05 ENE-D8 **S. STD. **
6	CS10 OMOFLUOROBENZENE **S. STD. **
7	C010 OMETHANE **
8	C015 OMETHANE
9	C020 L CHLORIDE *
10	C025 ROETHANE
11	C030 YLENE CHLORIDE
12	C251 LIN
13	C035 ONE
14	C252 LONITRILE
15	C040 ON DISULFIDE
16	C045 -DICHLOROETHENE *
17	C050 ICHLOROETHANE **
18	C055 S-1,2-DICHLOROETHENE
19	C000 RICHLOROFLUOROMETHANE
20	C060 ROFORM *
21	C065 -DICHLOROETHANE
22	C110 UTANONE
23	C115 -TRICHLOROETHANE
24	C120 ON TETRACHLORIDE
25	C125 L ACETATE
26	C130 D DICHLOROMETHANE
27	C140 -DICHLOROPROPANE *
28	C145 S-1,3 DICHLOROPROPENE
29	C150 HLOROETHENE
30	C155 ROMOCHLOROMETHANE
31	C160 2-TRICHLOROETHANE
32	C165 ENE
33	C143 -1,3-DICHLOROPROPENE
34	C175 -HLOROETHYL VINYL ETHER
35	C180 OFORM **
36	C220 TRACHLOROETHENE
37	C210 -EXANONE
38	C205 ETHYL 2-PENTANONE
39	C225 1,2,2-TETRACHLOROETHANE **
40	C230 LUENE *
41	C235 ROBENZENE **
42	C240 YL BENZENE *
43	C245 YRENE
44	C250 YLENES
45	C253 1,3-DICHLOROBENZENE
46	C254 1,4-DICHLOROBENZENE
47	C255 1,2-DICHLOROBENZENE

No	Name
48	C250 [illegible]/LENE
49	[illegible]ENE-D6 **S. STD. **

No	m/z	[illegible]	Time	Ref	RRT	Meth	Area(Hght)	Amount	%Tot
1	49	[illegible]	7:22	1	1.000	M XX	15702.	50.000 NG	2.08
2	114	[illegible]	9:38	2	1.000	A BB	113786.	50.000 NG	2.08
3	117	[illegible]	21:48	3	1.000	A BB	82666.	50.000 NG	2.08
4	65	[illegible]	8:36	1	1.168	A BB	22356.	50.000 NG	2.08
5	98	[illegible]	14:49	3	0.679	A BB	101320.	50.000 NG	2.08
6	95	[illegible]	28:33	3	1.309	A BB	77589.	50.000 NG	2.08
7	50	[illegible]	2:28	1	0.336	A BV	18209.	50.000 NG	2.08
8	94	[illegible]	3:00	1	0.408	M XX	49677.	50.000 NG	2.08
9	62	[illegible]	2:36	1	0.353	M XX	31195.	50.000 NG	2.08
10	64	[illegible]	3:05	1	0.418	M XX	25081.	50.000 NG	2.08
11	49	[illegible]	4:37	1	0.627	A BB	29568.	50.000 NG	2.08
12	56	[illegible]	3:50	1	0.521	M XX	4861.	50.000 NG	2.08
13	43	[illegible]	3:53	1	0.527	M XX	14594.	50.000 NG	2.08
14	53	[illegible]	4:51	1	0.658	A BB	6512.	50.000 NG	2.08
15	76	[illegible]	4:37	1	0.627	A BB	80436.	50.000 NG	2.08
16	96	[illegible]	4:01	1	0.545	A BB	7146.	50.000 NG	2.08
17	63	[illegible]	5:36	1	0.760	M XX	13495.	50.000 NG	2.08
18	61	[illegible]	5:00	1	0.678	A BB	24999.	50.000 NG	2.08
19	101	[illegible]	3:17	1	0.445	A BB	9542.	50.000 NG	2.08
20	83	[illegible]	6:58	1	0.945	A BB	13321.	50.000 NG	2.08
21	62	[illegible]	8:50	1	1.199	A BB	27652.	50.000 NG	2.08
22	43	[illegible]	6:23	2	0.662	A BB	13930.	50.000 NG	2.08
23	97	[illegible]	7:45	2	0.804	A BB	69949.	50.000 NG	2.08
24	117	[illegible]	8:18	2	0.861	A BB	58185.	50.000 NG	2.08
25	43	[illegible]	5:38	2	0.584	A BB	59862.	50.000 NG	2.08
26	83	[illegible]	11:53	2	1.233	A BB	68794.	50.000 NG	2.08
27	63	[illegible]	11:06	2	1.152	A BB	32854.	50.000 NG	2.08
28	75	[illegible]	14:02	2	1.455	A BB	47833.	50.000 NG	2.08
29	95	[illegible]	10:28	2	1.086	A BB	42644.	50.000 NG	2.08
30	129	[illegible]	19:02	2	1.974	A BB	51218.	50.000 NG	2.08
31	97	[illegible]	16:53	2	1.751	A BB	32070.	50.000 NG	2.08
32	78	[illegible]	8:47	2	0.911	A BB	91386.	50.000 NG	2.08
33	75	[illegible]	16:17	2	1.688	A BB	40226.	50.000 NG	2.08
34	63	[illegible]	13:18	2	1.380	M XX	2274.	50.000 NG	2.08
35	173	[illegible]	27:00	2	2.801	A BB	38042.	50.000 NG	2.08
36	166	[illegible]	17:52	3	0.819	A BB	46941.	50.000 NG	2.08
37	43	[illegible]	13:21	3	0.612	A BB	23400.	50.000 NG	2.08
38	43	[illegible]	17:13	3	0.789	A BB	17512.	50.000 NG	2.08
39	83	[illegible]	28:31	3	1.308	A BB	51776.	50.000 NG	2.08
40	91	[illegible]	15:09	3	0.694	A BB	105869.	50.000 NG	2.08
41	112	[illegible]	22:00	3	1.009	A BB	68042.	50.000 NG	2.08
42	106	[illegible]	22:25	3	1.028	A BB	29306.	50.000 NG	2.08
43	104	[illegible]	25:25	3	1.166	A BB	69393.	50.000 NG	2.08
44	106	[illegible]	22:49	3	1.046	A BB	87571.	50.000 NG	2.08
45	146	[illegible]	34:29	3	1.581	A BB	77948.	50.000 NG	2.08
46	146	[illegible]	35:05	3	1.609	A BB	79034.	50.000 NG	2.08
47	146	[illegible]	36:40	3	1.682	A BB	70481.	50.000 NG	2.08
48	106	[illegible]	25:05	3	1.150	A BB	44660.	50.000 NG	2.08
49	NOT FOUND								

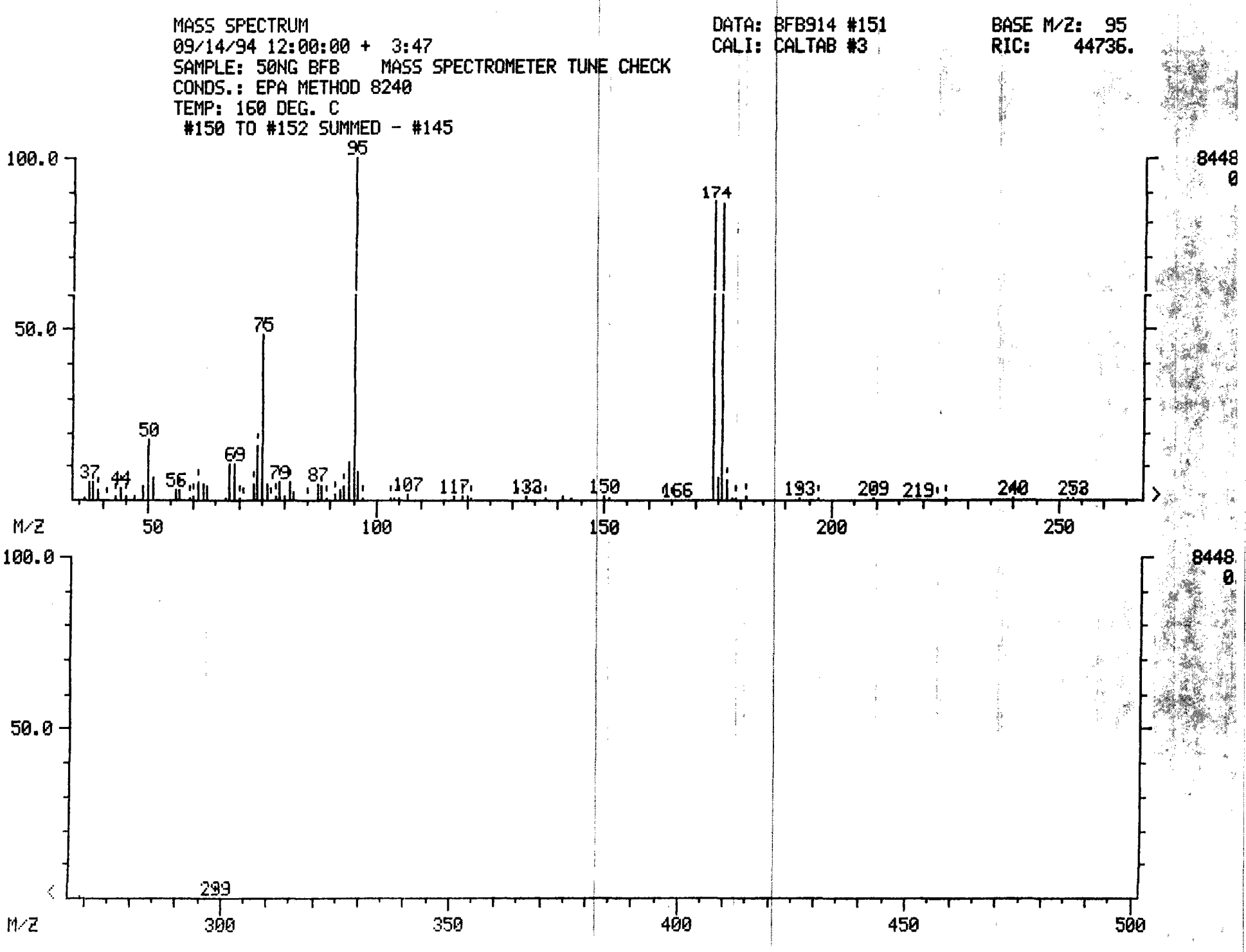
BROMOFLUOROBENZENE

Tuning Report
9/14/94 12:00:00 + 3:47
Instrument: FINN
#150 to #152 summed - #145
Case Number:

Data: BFB914 # 151
Cali: CALTAB # 3
Analyst: UC
Laboratory:

Base m/z: 95
RIC: 44736.
Acct. No.: GC
Contract:

			Ion Abundance Criteria				
m/z	Intensity	% RA	Min %	Max %	Mass	Actual	Status
50	1542.	18.3	15.0	40.0	95	18.3	PASS
75	4096.	48.5	30.0	60.0	95	48.5	PASS
95	8448.	100.0	100.0	---	---	100.0	PASS
96	697.	8.3	5.0	9.0	95	8.3	PASS
173	0.	0.0	---	2.0	174	0.0	PASS
174	7296.	86.4	50.0	---	95	86.4	PASS
175	542.	6.4	5.0	9.0	174	7.4	PASS
176	7248.	85.8	95.0	101.0	174	99.3	PASS
177	503.	6.0	5.0	9.0	176	6.9	PASS

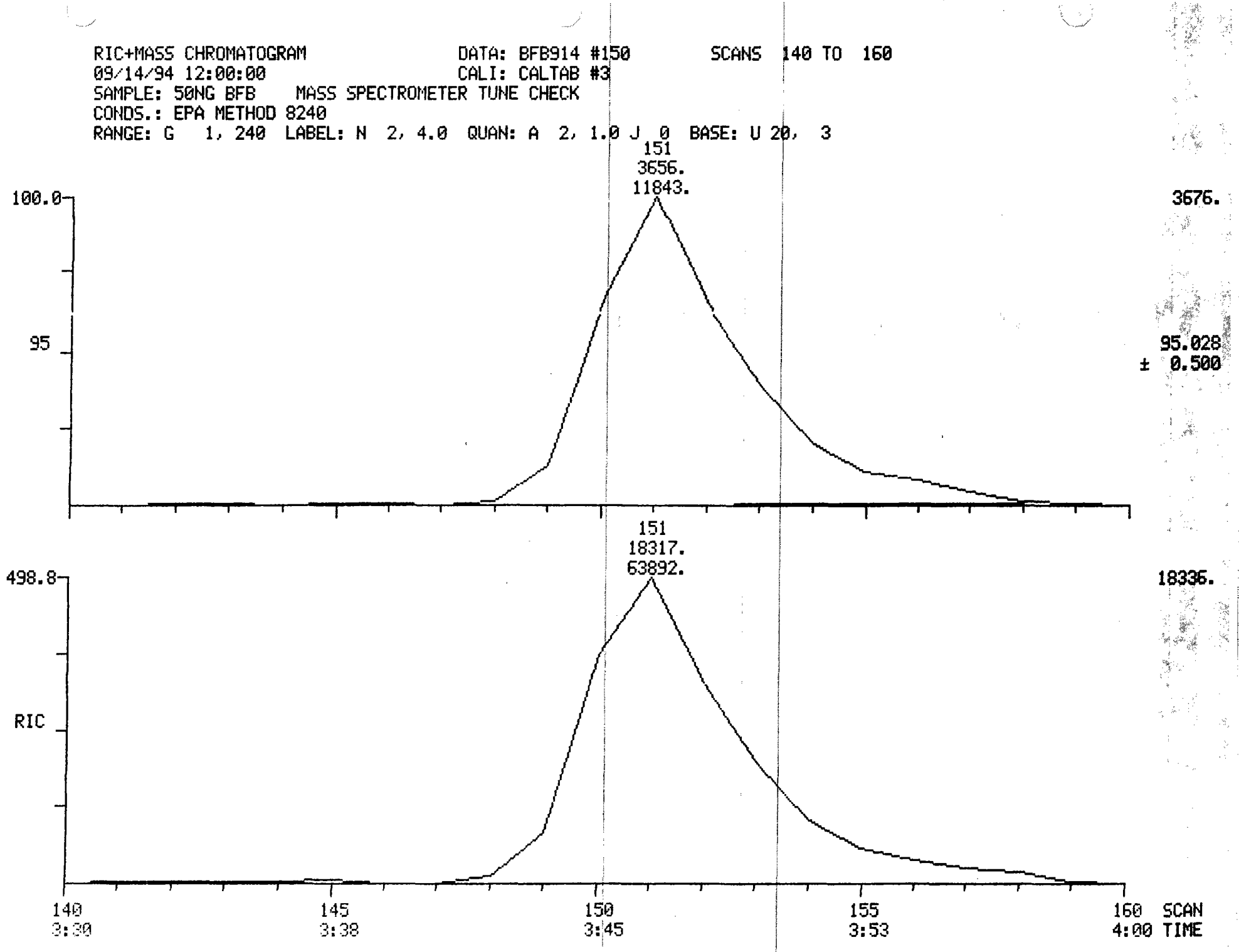

MASS SPECTRUM
09/14/94 12:00:00 + 3:47
SAMPLE: 50NG BFB MASS SPECTROMETER TUNE CHECK
CONDS.: EPA METHOD 8240
TEMP: 160 DEG. C
#150 TO #152 SUMMED - #145
DATA: BFB914 #151
CALI: CALTAB #3
BASE M/Z: 95
RIC: 44736.
100.0
50.0
95
174
75
50
69
37
44
56
79
87
107
117
133
150
166
193
209
219
240
253
8448
0
M/Z
50
100
150
200
250
100.0
50.0
8448
0
293
M/Z
300
350
400
450
500

Mass List
09/14/94 12:00:00 + 3:47
Sample: 50NG BFB
nds.: EPA METHOD 8240
#150 to #152 summed - #145

Data: BFB914 # 151
Cali: CALTAB # 3
MASS SPECTROMETER TUNE CHECK

Base m/z: 95
RIC: 44736.

36		0.00	0.	Minima	Min Inten:		0.
299				Maxima	# 0		
Mass		% RA	Inten.	Mass		% RA	Inten.
36?		0.37	31.	111		0.19	16.
37?		5.68	480.	117		1.11	94.
38?		5.35	452.	119	S	0.91	77.
39?	S	2.70	228.	120		0.91	77.
41?	S	0.25	21.	121	S	0.31	26.
43?	S	0.96	81.	122		0.22	19.
44?	S	3.53	298.	123		0.22	19.
45?	S	1.02	86.	131	S	0.25	21.
47?		1.02	86.	133	S	1.43	121.
48?		0.19	16.	135	S	0.27	23.
49?		4.21	356.	136	S	0.26	22.
50?		18.25	1542.	137	S	0.46	39.
51?		6.37	538.	141		1.20	101.
56?		3.29	278.	143		0.85	72.
57?	S	3.10	262.	150		0.91	77.
59?	S	0.76	64.	151		0.52	44.
60?	S	1.01	85.	163	S	0.09	8.
61?	S	5.66	478.	165	S	0.02	2.
62?		4.59	388.	166		0.18	15.
63?		4.06	343.	174		86.36	7296.
67?		0.60	51.	175		6.42	542.
68?		10.52	889.	176		85.80	7248.
69	S	10.89	920.	177	S	5.95	503.
70	S	0.62	52.	178		0.54	46.
71	S	0.21	18.	179	S	0.53	45.
73	S	4.92	416.	181	S	1.12	95.
74	S	16.29	1376.	193	S	0.51	43.
75	S	48.48	4096.	197	S	0.46	39.
76		4.64	392.	209	S	0.44	37.
77	S	0.30	25.	219	S	0.26	22.
78	S	0.91	77.	223	S	0.24	20.
79	S	5.23	442.	225	S	0.45	38.
80		1.16	98.	239	S	0.01	1.
81	S	5.15	435.	240	S	0.34	29.
82		2.21	187.	241	S	0.27	23.
85	S	0.09	8.	252		0.40	34.
87		4.69	396.	253	S	0.60	51.
88		3.89	329.	255	S	0.01	1.
89	S	0.33	28.	269		0.50	42.
91	S	1.57	133.	299	S	0.22	19.
92		3.03	256.				
93	S	4.38	370.				
94		11.35	959.				
95	S	100.00	8448.				
96		8.25	697.				
97	S	0.50	42.				
103	S	0.71	60.				
104		0.31	26.				
105	S	0.85	72.				
107		1.73	146.				


RIC+MASS CHROMATOGRAM
09/14/94 12:00:00
SAMPLE: 50NG BFB MASS SPECTROMETER TUNE CHECK
CONDS.: EPA METHOD 8240
RANGE: G 1, 240 LABEL: N 2, 4.0 QUAN: A 2, 1.0 J 0 BASE: U 20, 3
DATA: BFB914 #150
CALI: CALTAB #3
SCANS 140 TO 160
151
3656.
11843.
100.0
95
3676.
95.028
± 0.500
151
18317.
63892.
498.8
RIC
18336.
140
3:30
145
3:38
150
3:45
155
3:53
160 SCAN
4:00 TIME

BROMOFLUOROBENZENE

Tuning Report
[illegible]3/94 10:55:00 + 3:45
Instrument: FINN
#149 to #151 summed
Case Number:

Data: BFB923 # 150
Cali: CALTAB # 3
Analyst: UC
Laboratory:

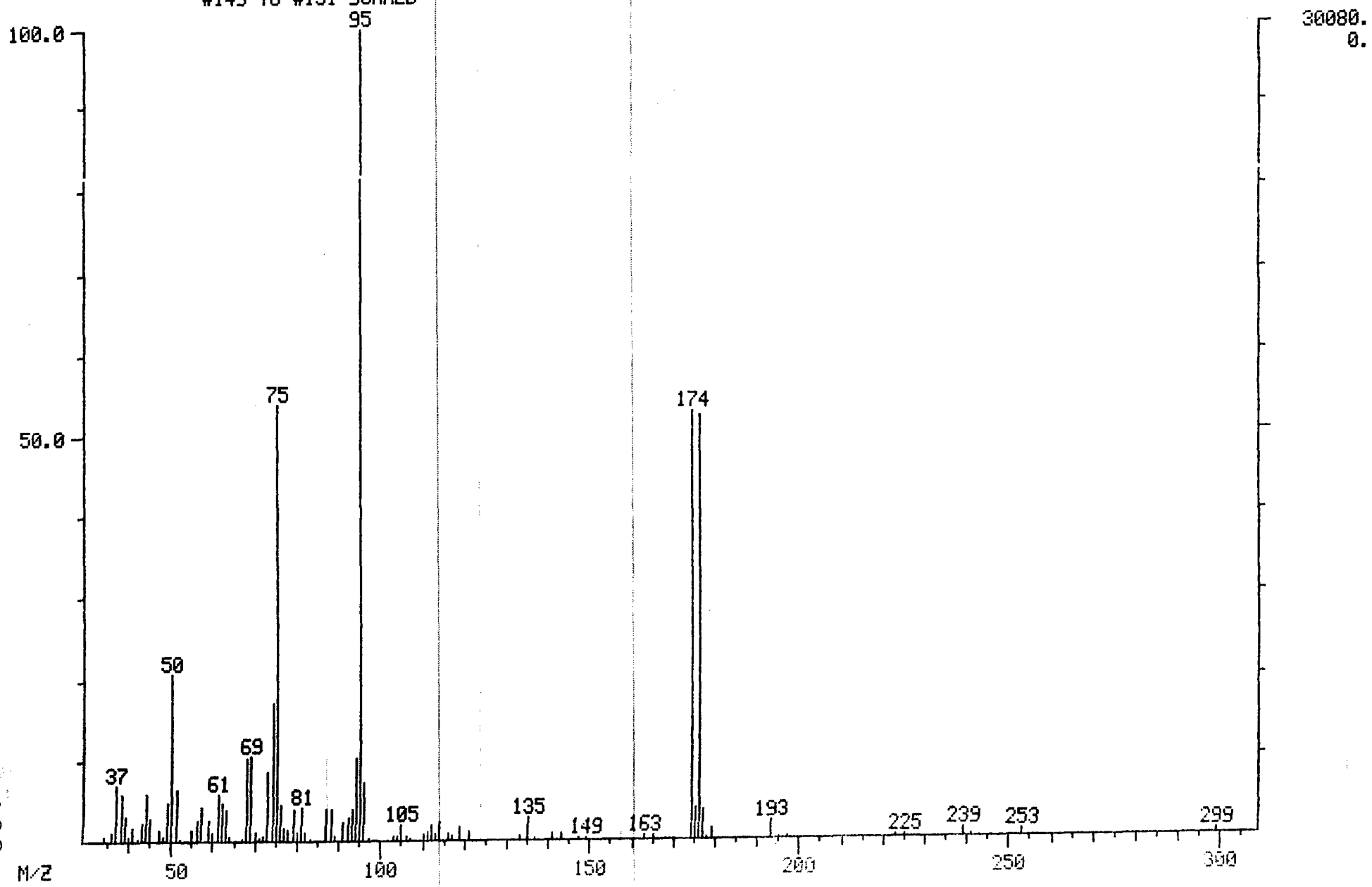
Base m/z: 95
RIC: 146176.
Acct. No.: 4667-001
Contract:

m/z	Intensity	% RA	Ion Abundance Criteria Min %	Max %	Mass	Actual	Status
50	6224.	20.7	15.0	40.0	95	20.7	PASS
75	16208.	53.9	30.0	60.0	95	53.9	PASS
95	30080.	100.0	100.0	---	---	100.0	PASS
96	2200.	7.3	5.0	9.0	95	7.3	PASS
173	0.	0.0	---	2.0	174	0.0	PASS
174	15856.	52.7	50.0	---	95	52.7	PASS
175	1144.	3.8	5.0	9.0	174	7.2	PASS
176	15744.	52.3	95.0	101.0	174	99.3	PASS
177	1100.	3.7	5.0	9.0	176	7.0	PASS

MASS SPECTRUM
09/23/94 10:55:00 + 3:45
SAMPLE: 50NG BFB MASS SPECTROMETER TUNE CHECK
CONDS.: EPA METHOD 8240
TEMP: 160 DEG. C
#149 TO #151 SUMMED

DATA: BFB923 #150
CALI: CALTAB #3

BASE M/Z: 95
RIC: 146176.



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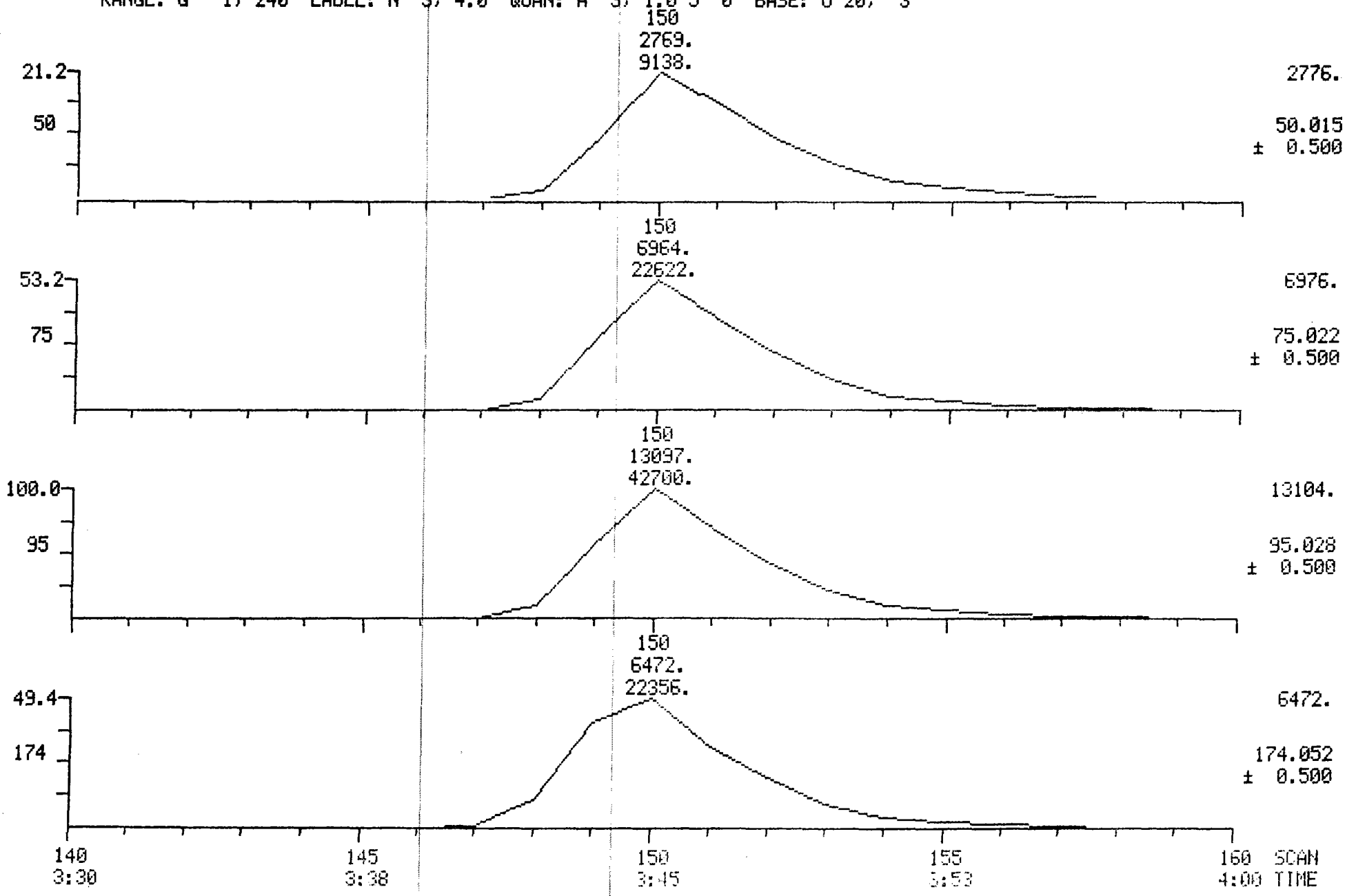
Mass List Data: BFB923 # 150 Base m/z: 95
09/23/94 10:55:00 + 3:45 Cali: CALTAB # 3 RIC: 146176.
Sample: 50NG BFB MASS SPECTROMETER TUNE CHECK
Comnds.: EPA METHOD 8240
#149 to #151 summed

34	0.00	0.	Minima	Min Inten:	0.
299			Maxima	# 0	
Mass	% RA	Inten.	Mass	% RA	Inten.
34?	0.60	179.	94	10.31	3100.
36?	1.22	368.	95	100.00	30080.
37?	6.87	2068.	96	7.31	2200.
38?	5.66	1704.	97	0.18	54.
39?	3.07	922.	103	0.58	175.
40?	0.45	136.	104	0.54	161.
41?	1.78	534.	105	1.96	589.
42?	0.27	82.	106	0.55	166.
43?	2.16	651.	107	0.39	116.
44?	5.74	1726.	110	0.88	265.
45?	2.77	833.	111	1.09	328.
47?	1.39	417.	112	1.90	572.
48?	0.44	131.	113	0.96	289.
49?	4.64	1396.	114	2.12	637.
50?	20.69	6224.	116	0.82	247.
51?	6.46	1944.	117	0.59	178.
52?	0.23	69.	118	0.13	40.
53?	0.05	15.	119	1.77	533.
55?	1.38	415.	121	1.21	363.
56?	2.44	734.	128	0.05	15.
57?	4.06	1222.	133	0.45	136.
59?	2.47	742.	135	2.69	809.
60?	1.24	374.	137	0.21	63.
61?	5.70	1714.	141	0.81	244.
62?	4.61	1388.	143	0.86	260.
63?	3.86	1162.	145	0.13	38.
64?	0.64	194.	147	0.15	46.
67?	0.26	78.	149	0.25	74.
68?	10.17	3060.	163	0.64	192.
69	10.45	3144.	165	0.56	167.
70	1.06	320.	174	52.71	15856.
71	0.25	76.	175	3.80	1144.
72	0.52	155.	176	52.34	15744.
73	8.70	2616.	177	3.66	1100.
74	17.21	5176.	178	0.15	45.
75	53.88	16208.	179	1.38	415.
76	4.53	1364.	180	0.05	15.
77	1.55	466.	181	0.08	23.
78	1.26	380.	193	2.19	660.
79	3.95	1188.	194	0.13	40.
80	1.18	354.	195	0.31	94.
81	4.13	1242.	197	0.25	74.
82	1.10	330.	223	0.26	78.
83	0.20	61.	225	0.35	104.
87	3.86	1162.	237	0.05	16.
88	3.91	1176.	239	1.20	360.
89	0.61	182.	241	0.25	74.
91	2.26	680.	253	0.95	287.
92	2.70	811.	299	0.56	167.
93	3.95	1188.			

MASS CHROMATOGRAMS DATA: BFB923 #149 SCANS 140 TO 160
09/23/94 10:55:00 CALI: CALTAB #3
SAMPLE: 50NG BFB MASS SPECTROMETER TUNE CHECK
CONDS.: EPA METHOD 8240
RANGE: G 1, 240 LABEL: N 3, 4.0 QUAN: A 3, 1.0 J 0 BASE: U 20, 3



BROMOFLUOROBENZENE

)ning Report
09/29/94 8:45:00 + 3:44
Instrument: FINN
#148 to #150 summed
Case Number:

Data: BFB929 # 149
Cali: CALTAB # 3
Analyst: UC
Laboratory:

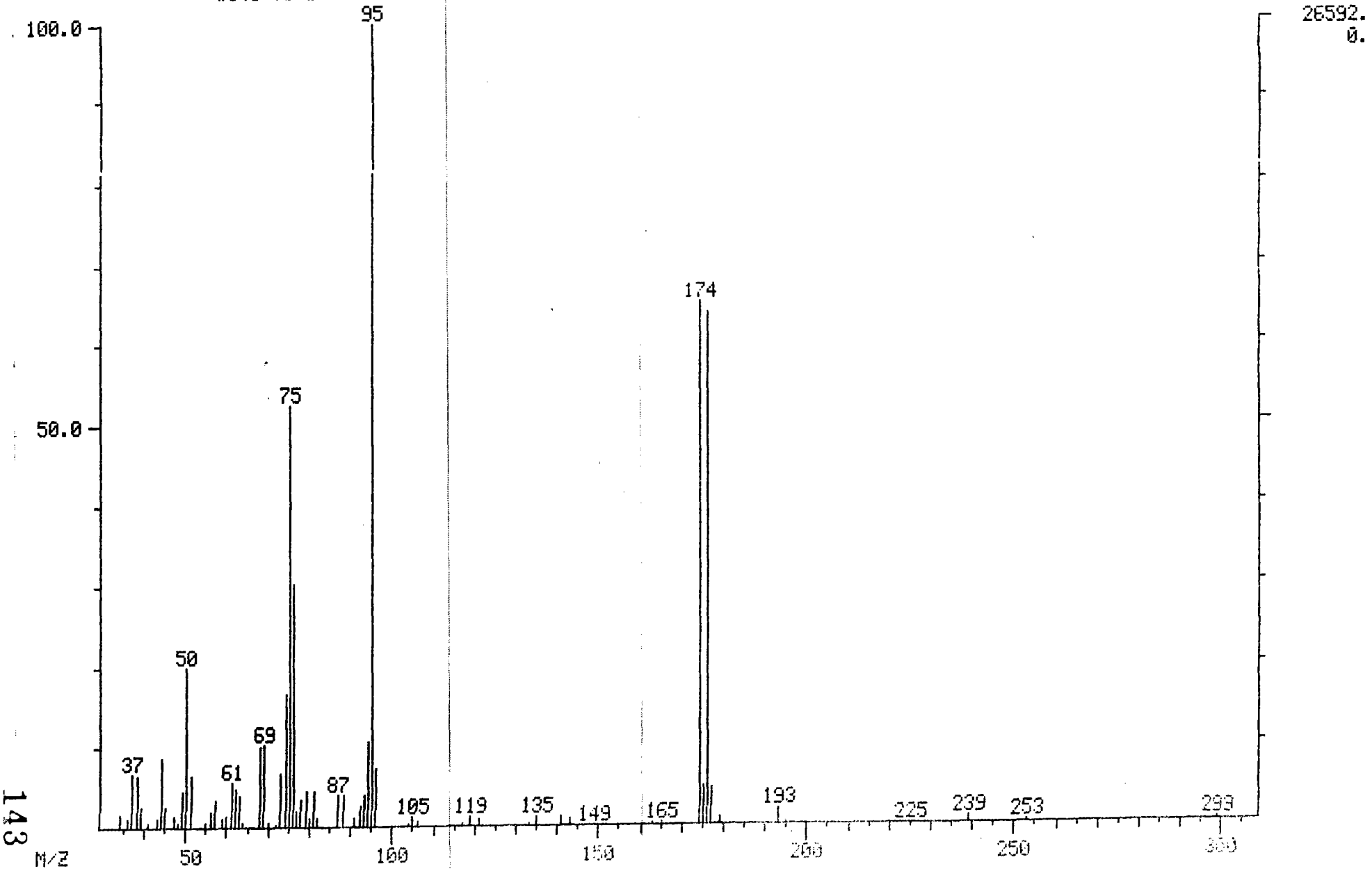
Base m/z: 95
RIC: 136448.
Acct. No.: INTERNAL
Contract:

m/z	Intensity	% RA	Ion Abundance Criteria Min %	Max %	Mass	Actual	Status
50	5304.	19.9	15.0	40.0	95	19.9	PASS
75	13936.	52.4	30.0	60.0	95	52.4	PASS
95	26592.	100.0	100.0	---	---	100.0	PASS
96	1876.	7.1	5.0	9.0	95	7.1	PASS
173	0.	0.0	---	2.0	174	0.0	PASS
174	17344.	65.2	50.0	---	95	65.2	PASS
175	1234.	4.6	5.0	9.0	174	7.1	PASS
176	16960.	63.8	95.0	101.0	174	97.8	PASS
177	1146.	4.3	5.0	9.0	176	6.8	PASS

MASS SPECTRUM
09/29/94 8:45:00 + 3:44
SAMPLE: 50NG BFB MASS SPECTROMETER TUNE CHECK
CONDS.: EPA METHOD 624
TEMP: 160 DEG. C
#148 TO #150 SUMMED

DATA: BFB929 #149
CALI: CALTAB #3

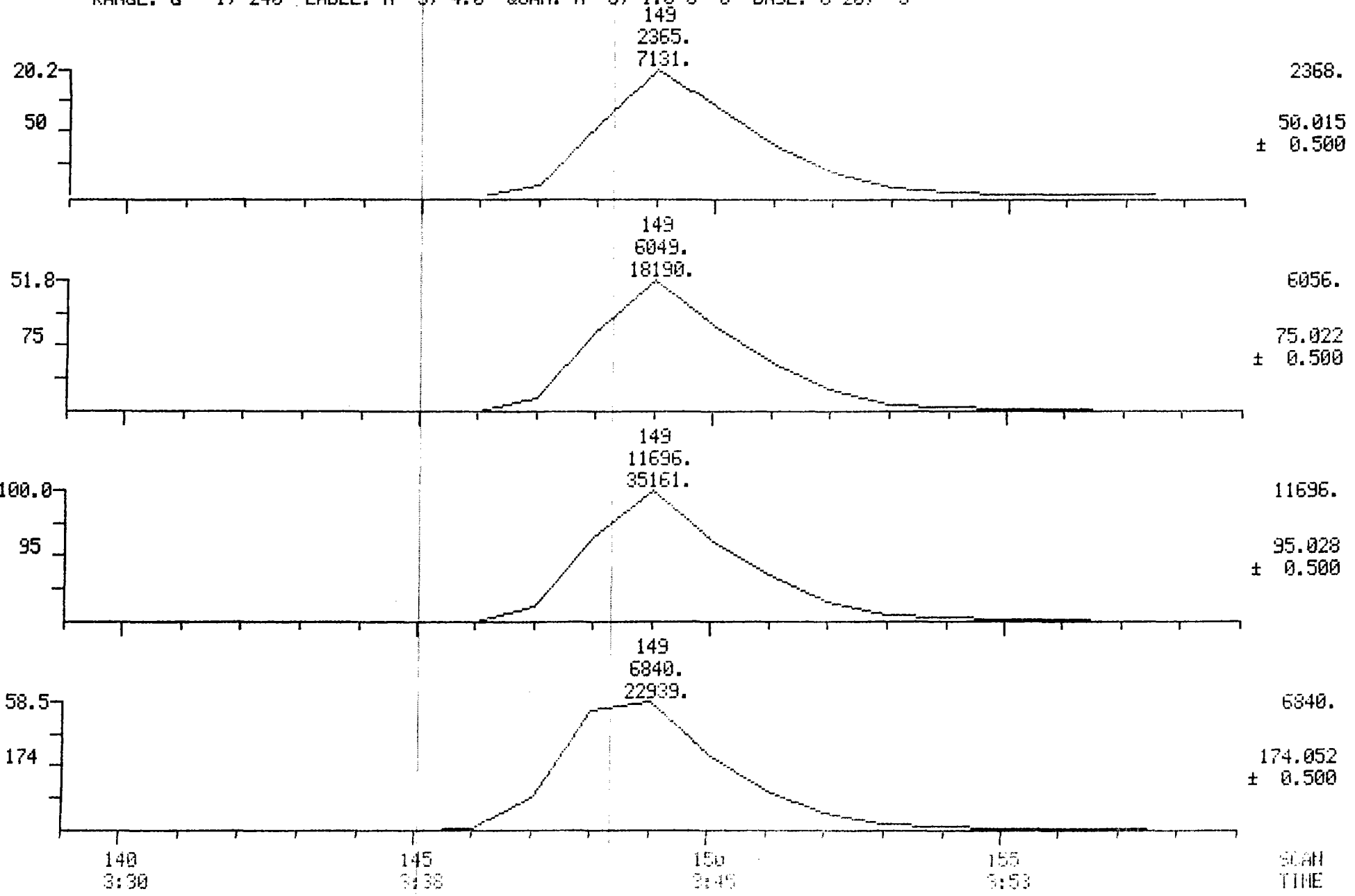
BASE M/Z: 95
RIC: 136448.



Mass List Data: BFB929 # 149 Base m/z: 95
9/29/94 8:45:00 + 3:44 Cali: CALTAB # 3 RIC: 136448.
Sample: 50NG BFB MASS SPECTROMETER TUNE CHECK
Cmds.: EPA METHOD 624
#148 to #150 summed

34	0.00	0.	Minima	Min Inten:	0.
299			Maxima	# 0	
Mass	% RA	Inten.	Mass	% RA	Inten.
34?	1.65	440.	117	0.36	95.
36?	1.14	304.	118	0.12	32.
37?	6.57	1746.	119	1.24	330.
38?	6.26	1664.	121	0.72	192.
39?	2.47	657.	133	0.14	37.
40?	0.07	19.	135	1.08	286.
41?	0.46	121.	141	1.00	266.
43?	0.99	263.	143	0.88	235.
44?	8.47	2252.	149	0.06	17.
45?	2.35	626.	163	0.15	41.
47?	1.50	399.	165	0.34	91.
48?	0.51	135.	174	65.22	17344.
49?	4.52	1202.	175	4.64	1234.
50?	19.95	5304.	176	63.78	16960.
51?	6.27	1668.	177	4.31	1146.
55?	0.61	162.	179	0.94	249.
56?	1.80	479.	193	1.85	491.
57?	3.29	875.	195	0.19	51.
59?	1.20	320.	225	0.07	19.
60?	1.33	353.	239	0.95	252.
61?	5.40	1436.	253	0.33	87.
62?	4.82	1282.	299	0.15	41.
63?	3.90	1036.			
64?	0.64	169.			
67?	0.06	17.			
68?	9.94	2644.			
69	10.35	2752.			
70	0.63	167.			
72	0.33	89.			
73	6.51	1732.			
74	16.70	4440.			
75	52.41	13936.			
76	30.26	8048.			
77	1.82	483.			
78	3.26	867.			
79	4.53	1204.			
80	1.23	328.			
81	4.38	1166.			
82	1.06	283.			
87	3.90	1036.			
88	3.89	1034.			
91	1.24	329.			
92	2.54	675.			
93	4.00	1064.			
94	10.57	2812.			
95	100.00	26592.			
96	7.05	1876.			
104	0.19	51.			
105	0.98	261.			
106	0.42	111.			

MASS CHROMATOGRAMS DATA: BFB929 #148 SCANS 139 TO 159
09/29/94 8:45:00 CALI: CALTAB #3
SAMPLE: 50NG BFB MASS SPECTROMETER TUNE CHECK
CONDS.: EPA METHOD 624
RANGE: G 1, 240 LABEL: N 3, 4.0 QUAN: A 3, 1.0 J 0 BASE: U 20, 3



BROMOFLUOROBENZENE

ing Report
4/94 11:42:00 + 3:43
trument: FINN
146 to #148 summed
ase Number:

Data: BFB1004B # 147
Cali: CALTAB # 3
Analyst: KE
Laboratory:

Base m/z: 95
RIC: 128512.
Acct. No.: -
Contract:

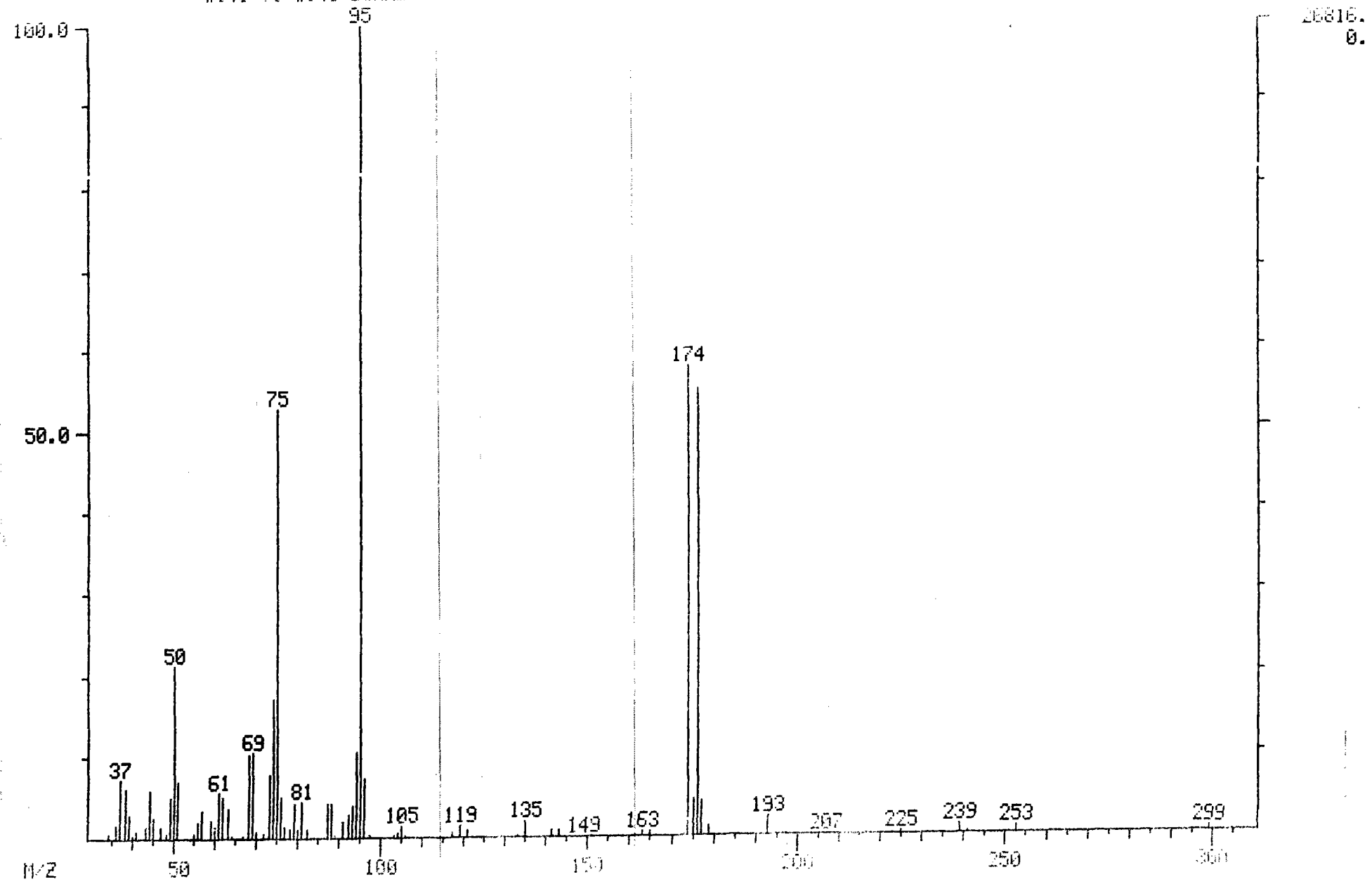
Ion Abundance Criteria

m/z	Intensity	% RA	Min %	Max %	Mass	Actual	Status
50	5704.	21.3	15.0	40.0	95	21.3	PASS
75	14128.	52.7	30.0	60.0	95	52.7	PASS
95	26816.	100.0	100.0	---	---	100.0	PASS
96	1948.	7.3	5.0	9.0	95	7.3	PASS
173	0.	0.0	---	2.0	174	0.0	PASS
174	15472.	57.7	50.0	---	95	57.7	PASS
175	1174.	4.4	5.0	9.0	174	7.6	PASS
176	14752.	55.0	95.0	101.0	174	95.3	PASS
177	1088.	4.1	5.0	9.0	176	7.4	PASS

MASS SPECTRUM
10/04/94 11:42:00 + 3:43
SAMPLE: 50NG BFB MASS SPECTROMETER TUNE CHECK
CONDS.: EPA METHOD 8240
TEMP: 160 DEG. C
#145 TO #148 SUMMED

DATA: BFB1004B #147
CALI: CALTAB #3

BASE M/Z: 95
RIC: 128512.



147

Mass List
10/04/94 11:42:00 + 3:43
Sample: 50NG BFB
Anas.: EPA METHOD 8240
#1-6 to #148 summed

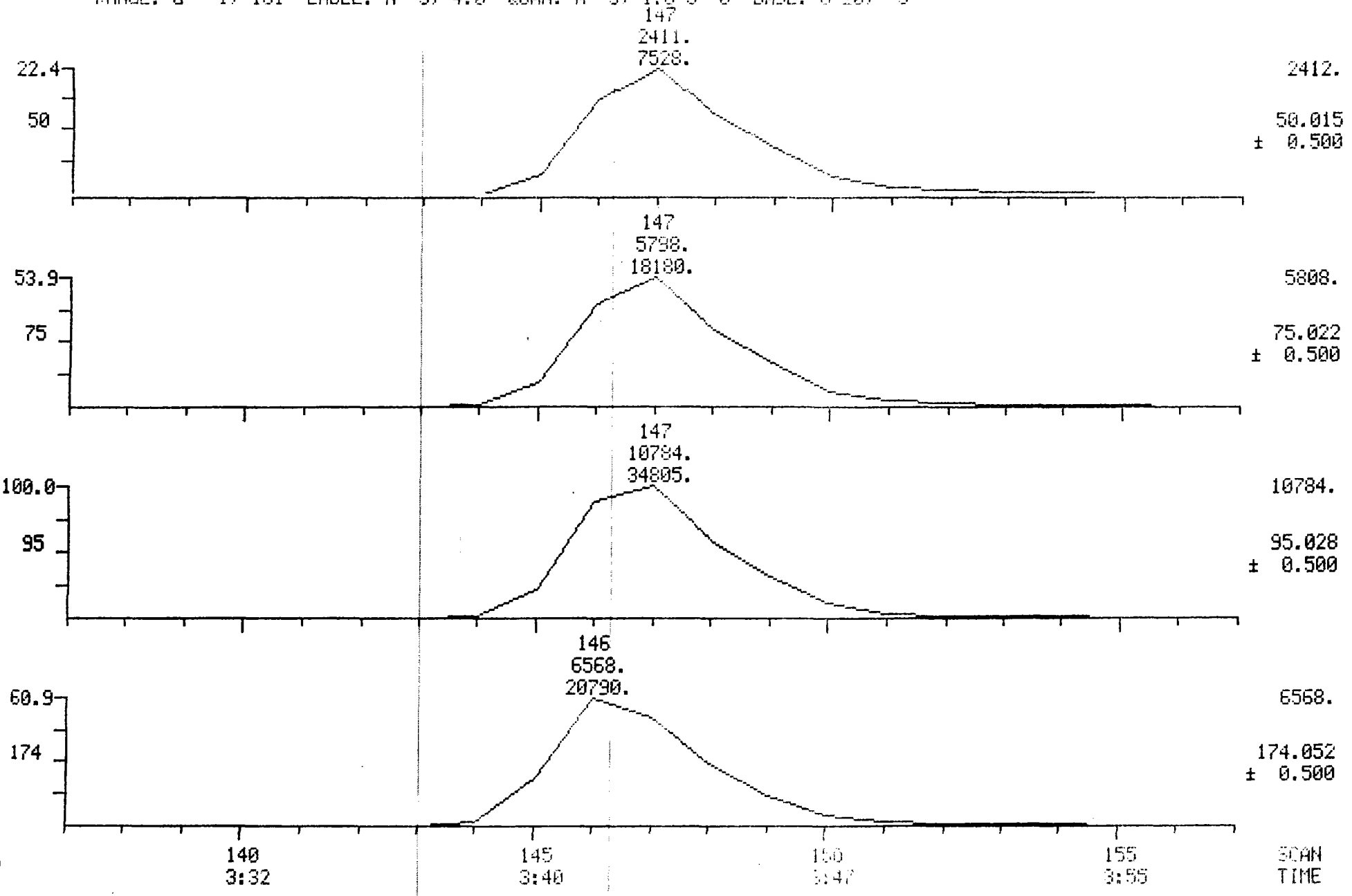
Data: BFB1004B # 147
Cali: CALTAB # 3

MASS SPECTROMETER TUNE CHECK

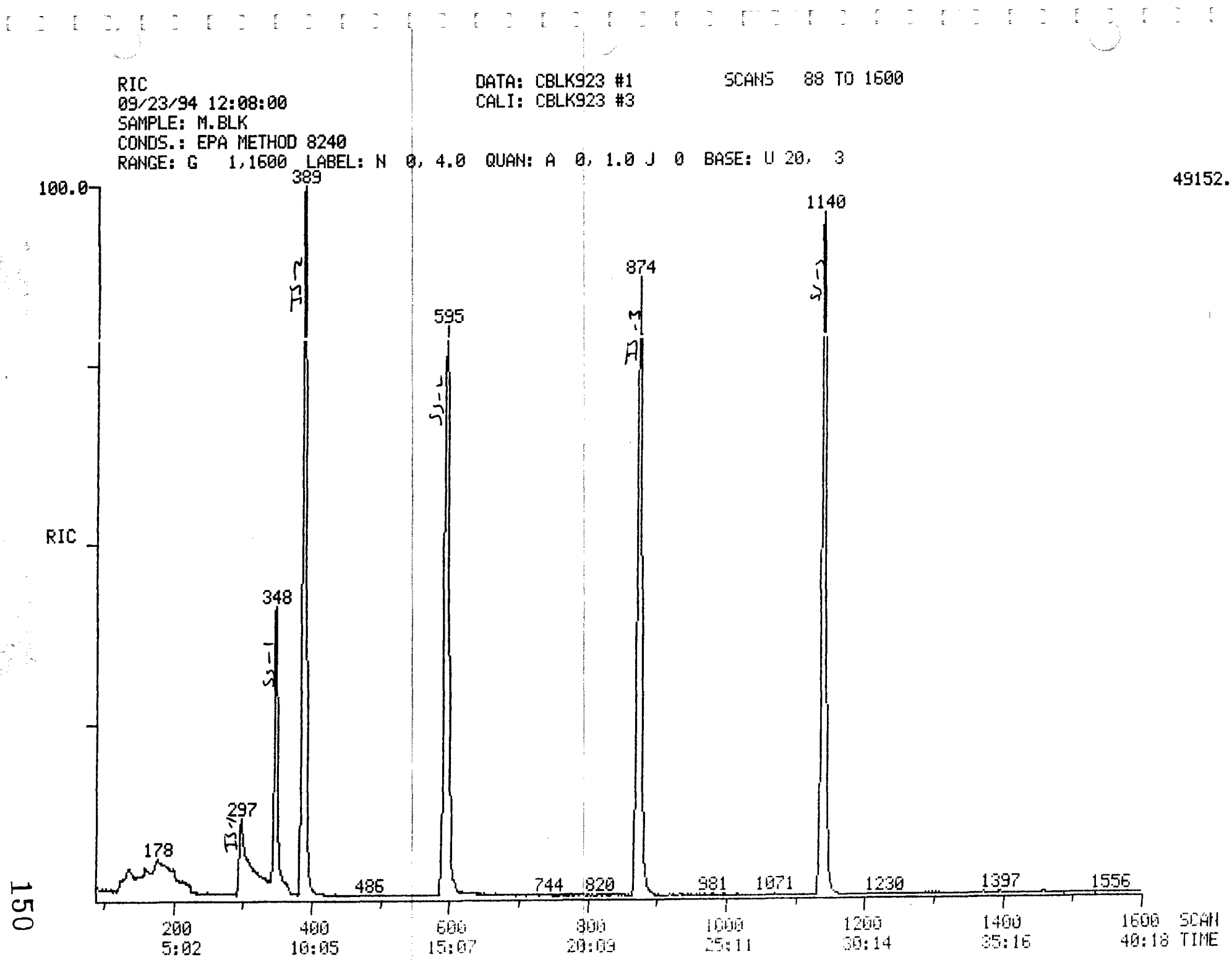
Base m/z: 95
RIC: 128512.

34	0.00	0.	Minima	Min Inten:	0.
300			Maxima #	0	
Mass	% RA	Inten.	Mass	% RA	Inten.
34?	0.59	157.	96	7.26	1948.
36?	1.67	449.	97	0.16	43.
37?	7.09	1902.	103	0.29	77.
38?	6.08	1630.	104	0.43	116.
39?	2.80	752.	105	1.42	381.
40?	0.25	68.	106	0.14	38.
41?	0.81	216.	107	0.10	27.
42?	0.08	22.	115	0.06	16.
43?	1.42	381.	117	0.59	157.
44?	5.91	1584.	118	0.06	15.
45?	2.56	687.	119	1.33	356.
47?	1.48	398.	120	0.08	22.
48?	0.61	163.	121	0.74	198.
49?	5.09	1366.	128	0.08	22.
50?	21.27	5704.	133	0.34	90.
51?	6.88	1846.	135	2.03	544.
52?	0.13	36.	137	0.09	24.
55?	0.65	175.	141	0.72	193.
56?	1.93	518.	143	0.80	214.
57?	3.29	882.	149	0.07	19.
59?	2.13	570.	163	0.54	144.
60?	1.25	334.	165	0.47	125.
61?	5.57	1494.	174	57.70	15472.
62?	5.09	1366.	175	4.38	1174.
63?	3.72	997.	176	55.01	14752.
64?	0.25	66.	177	4.06	1088.
67?	0.21	57.	178	0.07	20.
68?	10.34	2772.	179	1.07	286.
69	10.37	2780.	181	0.12	31.
70	0.75	200.	193	2.14	573.
71	0.07	18.	194	0.13	34.
72	0.44	117.	195	0.14	37.
73	7.74	2076.	197	0.06	15.
74	17.24	4624.	207	0.09	23.
75	52.68	14128.	223	0.08	22.
76	4.95	1328.	225	0.21	55.
77	1.38	369.	237	0.06	15.
78	1.19	320.	239	1.06	285.
79	4.18	1122.	241	0.22	59.
80	1.23	330.	253	0.74	199.
81	4.29	1150.	255	0.10	26.
82	1.22	326.	299	0.61	164.
87	4.12	1104.	300	0.06	17.
88	4.16	1116.			
89	0.38	101.			
91	1.86	500.			
92	2.73	732.			
93	3.90	1046.			
94	10.41	2792.			
95	100.00	26816.			

MASS CHROMATOGRAMS DATA: BFB1004B #146 SCANS 137 TO 157
10/04/94 11:42:00 CALI: CALTAB #3
SAMPLE: 50NG BFB MASS SPECTROMETER TUNE CHECK
CONDS.: EPA METHOD 8240
RANGE: G 1, 181 LABEL: N 3, 4.0 QUAN: A 3, 1.0 J 0 BASE: U 30, 3



149


RIC
09/23/94 12:08:00
SAMPLE: M.BLK
CONDS.: EPA METHOD 8240
RANGE: G 1,1600 LABEL: N 0, 4.0 QUAN: A 0, 1.0 J 0 BASE: U 20, 3
DATA: CBLK923 #1
CALI: CBLK923 #3
SCANS 88 TO 1600
49152.
100.0
RIC
389
348
297
178
595
874
1140
486
744
820
981
1071
1230
1397
1556
200
5:02
400
10:05
600
15:07
800
20:09
1000
25:11
1200
30:14
1400
35:16
1600 SCAN
40:18 TIME
150

Quantitation Report File: CBLK923

Data: CBLK923.TI
?/23/94 12:08:00
Sample: M.BLK
Conds.: EPA METHOD 8240
Formula: 5ML Instrument: FINN Weight: 0.000
Submitted by: PTL Analyst: UC Acct. No.:

AMOUNT=AREA * REF AMNT/(REF AREA * RESP FACT)
Resp. fac. from Library Entry

No	Name
1	CI01 BROMOCHLOROMETHANE **INT. STD. **
2	CI10 1,4-DIFLUOROBENZENE **INT. STD. **
3	CI20 CHLOROBENZENE-D5 **INT. STD. **
4	CS15 1,2-DICHLOROETHANE-D4 **S. STD. **
5	CS05 TOLUENE-D8 **S. STD. **
6	CS10 4-BROMOFLUOROBENZENE **S. STD. **
7	C010 CHLOROMETHANE **
8	C015 BROMOMETHANE
9	C020 VINYL CHLORIDE *
10	C025 CHLOROETHANE
11	C030 METHYLENE CHLORIDE
12	C251 ACROLIN
13	C035 ACETONE
14	C252 ACRYLONITRILE
15	C040 CARBON DISULFIDE
16	C045 1,1-DICHLOROETHENE *
17	C050 1,1-DICHLOROETHANE **
18	C055 TRANS-1,2-DICHLOROETHENE
19	C000 TRICHLOROFLUOROMETHANE
20	C060 CHLOROFORM *
21	C065 1,2-DICHLOROETHANE
22	C110 2-BUTANONE
23	C115 1,1,1-TRICHLOROETHANE
24	C120 CARBON TETRACHLORIDE
25	C125 VINYL ACETATE
26	C130 BROMO DICHLOROMETHANE
27	C140 1,2-DICHLOROPROPANE *
28	C145 TRANS-1,3 DICHLOROPROPENE
29	C150 TRICHLOROETHENE
30	C155 DIBROMOCHLOROMETHANE
31	C160 1,1,2-TRICHLOROETHANE
32	C165 BENZENE
33	C143 CIS-1,3-DICHLOROPROPENE
34	C175 2-CHLOROETHYL VINYL ETHER
35	C180 BROMOFORM **
36	C220 TETRACHLOROETHENE
37	C210 2-HEXANONE
38	C205 4-METHYL 2-PENTANONE
39	C225 1,1,2,2-TETRACHLOROETHANE **
40	C230 TOLUENE *
41	C235 CHLOROBENZENE **
42	C240 ETHYL BENZENE *
43	C245 STYRENE
44	C250 M+P-XYLENES
45	C253 1,3-DICHLOROBENZENE
46	C254 1,4-DICHLOROBENZENE
47	C255 1,2-DICHLOROBENZENE

No	Name
48	C250 O-XYLENE
49	BENZENE-D6 **S. STD. **

No	m/z	Scan	Time	Ref	RRT	Meth	Area(Hght)	Amount	%Tot	
1	49	297	7:29	1	1.000	M XX	12703.	50.000 NG	16.18	
2	114	389	9:48	2	1.000	A BB	121792.	50.000 NG	16.18	
3	117	874	22:01	3	1.000	A BB	70127.	50.000 NG	16.18	
4	65	348	8:46	1	1.172	A BB	35778.	55.708 NG	18.03	111
5	98	595	14:59	3	0.681	A BB	97779.	52.086 NG	16.86	104
6	95	1140	28:43	3	1.304	M XX	72910.	50.265 NG	16.27	101
7	NOT FOUND									
8	NOT FOUND									
9	NOT FOUND									
10	NOT FOUND									
11	NOT FOUND									
12	NOT FOUND									
13	NOT FOUND									
14	NOT FOUND									
15	NOT FOUND									
16	NOT FOUND									
17	NOT FOUND									
18	NOT FOUND									
19	NOT FOUND									
20	NOT FOUND									
21	NOT FOUND									
22	NOT FOUND									
23	NOT FOUND									
24	NOT FOUND									
25	43	223	5:37	2	0.573	A VB	1017.	0.605 NG	0.20	
26	NOT FOUND									
27	NOT FOUND									
28	NOT FOUND									
29	NOT FOUND									
30	NOT FOUND									
31	NOT FOUND									
32	NOT FOUND									
33	NOT FOUND									
34	NOT FOUND									
35	NOT FOUND									
36	NOT FOUND									
37	NOT FOUND									
38	NOT FOUND									
39	NOT FOUND									
40	91	608	15:19	3	0.696	A BB	851.	0.414 NG	0.13	
41	NOT FOUND									
42	NOT FOUND									
43	NOT FOUND									
44	NOT FOUND									
45	NOT FOUND									
46	NOT FOUND									
47	NOT FOUND									
48	NOT FOUND									
49	NOT FOUND									

PROCEDURE: FILTER/TIC DIAGNOSTIC REPORT 9/23/94 12:59:48

DATA FILE: CBLK923

FILTER SCAN PARAMETERS

MAX. NUMBER TICS:	15
11-TABLE ENTRIES:	539
SCAN TOLERANCE :	2
MIN. RIC HT. [%]:	10
FIRST SCAN :	1
LAST SCAN :	1600
TIC THRESHOLD :	600

METHOD LIBRARY & LISTS

TIC I.S. LIBRARY:	LIBRARYLS
NBS SEARCH PROC :	SERLIB
PEAK FINDER PROC:	VOME
TCA I.S. LL :	LS
FILE NAME LIST :	TCAREF2

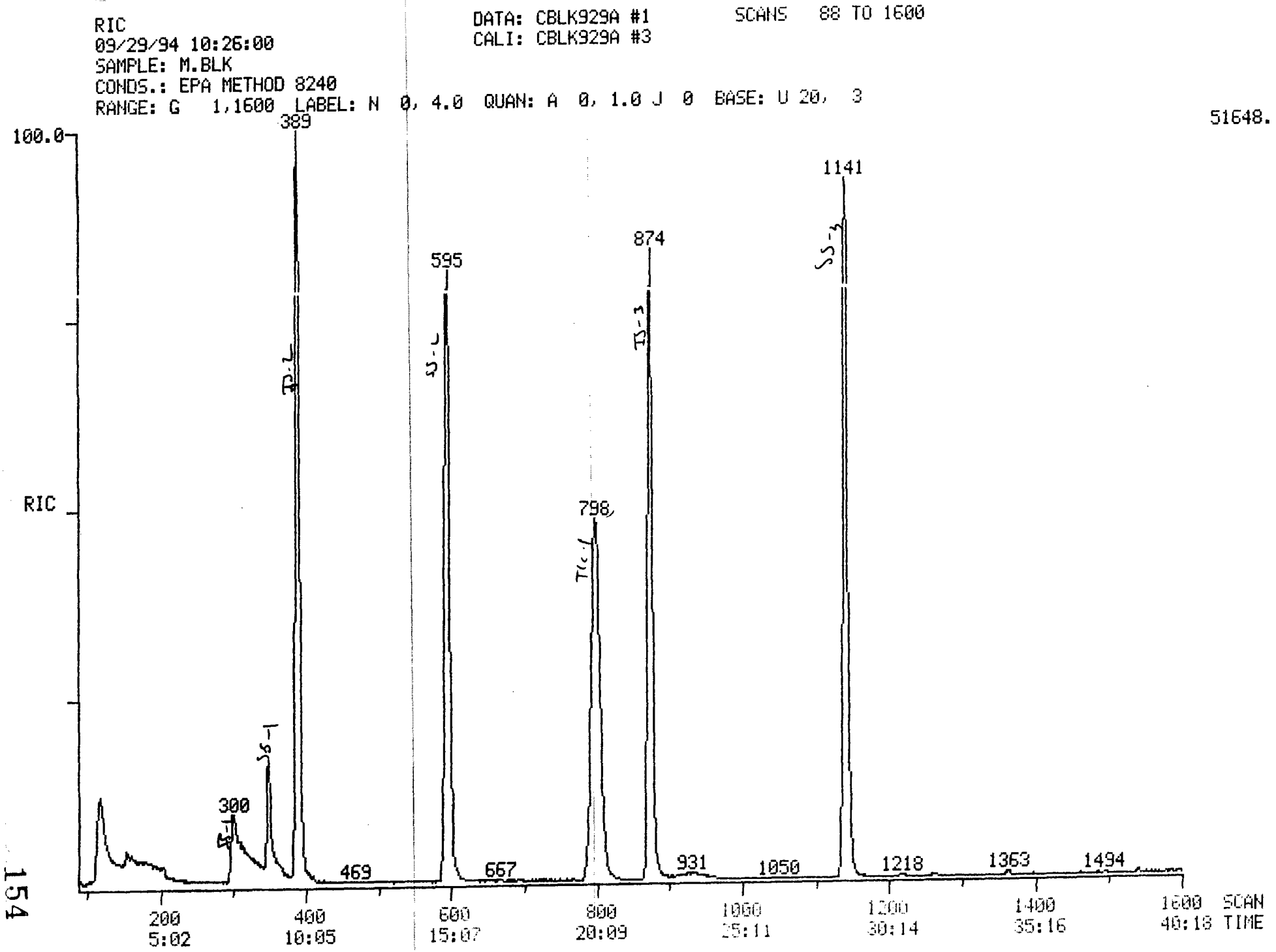
TARGET COMPOUND ANALYSIS:

TARGETS (QUAN LIST)	IS PEAKS	TOTAL TARGET PEAKS
5	3	8

FILTER PROCESSING:

	<-----REJECT PEAKS----->						
TOTAL PEAKS	< 1ST SCAN	> LAST SCAN	< MIN RIC HT	< SCAN TOL	> MAX # PEAKS	TOTAL REJECTS	TOTAL TICS
6	0	0	0	6	0	6	0

NO UNKNOWN PEAKS TO BE IDENTIFIED.


RIC
09/29/94 10:26:00
SAMPLE: M.BLK
CONDS.: EPA METHOD 8240
RANGE: G 1,1600 LABEL: N 0, 4.0 QUAN: A 0, 1.0 J 0 BASE: U 20, 3
DATA: CBLK929A #1
CALI: CBLK929A #3
SCANS 88 TO 1600
51648.
100.0
RIC
389
595
798
874
1141
300
469
667
931
1050
1218
1363
1494
200
5:02
400
10:05
600
15:07
800
20:09
1000
25:11
1200
30:14
1400
35:16
1600 SCAN
40:18 TIME

Quantitation Report File: CBLK929A

Data: CBLK929A.TI
9/29/94 10:26:00
Sample: M.BLK
Bonds.: EPA METHOD 8240
Formula: 5ML Instrument: FINN Weight: 0.000
Submitted by: PTL Analyst: UC Acct. No.:

AMOUNT=AREA * REF AMNT/(REF AREA * RESP FACT)
Resp. fac. from Library Entry

No	Name
1	CI01 BROMOCHLOROMETHANE **INT. STD. **
2	CI10 1,4-DIFLUOROBENZENE **INT. STD. **
3	CI20 CHLOROBENZENE-D5 **INT. STD. **
4	CS15 1,2-DICHLOROETHANE-D4 **S. STD. **
5	CS05 TOLUENE-D8 **S. STD. **
6	CS10 4-BROMOFLUOROBENZENE **S. STD. **
7	C010 CHLOROMETHANE **
8	C015 BROMOMETHANE
9	C020 VINYL CHLORIDE *
10	C025 CHLOROETHANE
11	C030 METHYLENE CHLORIDE
12	C251 ACROLIN
13	C035 ACETONE
14	C252 ACRYLONITRILE
15	C040 CARBON DISULFIDE
16	C045 1,1-DICHLOROETHENE *
17	C050 1,1-DICHLOROETHANE **
18	C055 TRANS-1,2-DICHLOROETHENE
19	C000 TRICHLOROFLUOROMETHANE
20	C060 CHLOROFORM *
21	C065 1,2-DICHLOROETHANE
22	C110 2-BUTANONE
23	C115 1,1,1-TRICHLOROETHANE
24	C120 CARBON TETRACHLORIDE
25	C125 VINYL ACETATE
26	C130 BROMO DICHLOROMETHANE
27	C140 1,2-DICHLOROPROPANE *
28	C145 TRANS-1,3 DICHLOROPROPENE
29	C150 TRICHLOROETHENE
30	C155 DIBROMOCHLOROMETHANE
31	C160 1,1,2-TRICHLOROETHANE
32	C165 BENZENE
33	C143 CIS-1,3-DICHLOROPROPENE
34	C175 2-CHLOROETHYL VINYL ETHER
35	C180 BROMOFORM **
36	C220 TETRACHLOROETHENE
37	C210 2-HEXANONE
38	C205 4-METHYL 2-PENTANONE
39	C225 1,1,2,2-TETRACHLOROETHANE **
40	C230 TOLUENE *
41	C235 CHLOROBENZENE **
42	C240 ETHYL BENZENE *
43	C245 STYRENE
44	C250 M+P-XYLENES
45	C253 1,3-DICHLOROBENZENE
46	C254 1,4-DICHLOROBENZENE
47	C255 1,2-DICHLOROBENZENE

Aug 29 A

No	Name
48	C250 O-XYLENE
49	BENZENE-D6 **S. STD. **

No	m/z	Scan	Time	Ref	RRT	Meth	Area(Hght)	Amount	%Tot	
1	49	300	7:33	1	1.000	A BB	10486.	50.000 NG	17.49	
2	114	389	9:48	2	1.000	A BB	121433.	50.000 NG	17.49	
3	117	873	21:59	3	1.000	A BB	89390.	50.000 NG	17.49	
4	65	348	8:46	1	1.160	A BB	16010.	45.142 NG	15.79	90
5	98	595	14:59	3	0.682	A BB	113897.	47.543 NG	16.63	93
6	95	1140	28:43	3	1.306	A BB	74554.	43.069 NG	15.06	96
7	NOT FOUND									
8	NOT FOUND									
9	NOT FOUND									
10	NOT FOUND									
11	NOT FOUND									
12	NOT FOUND									
13	NOT FOUND									
14	NOT FOUND									
15	NOT FOUND									
16	NOT FOUND									
17	NOT FOUND									
18	NOT FOUND									
19	NOT FOUND									
20	NOT FOUND									
21	NOT FOUND									
22	NOT FOUND									
23	NOT FOUND									
24	NOT FOUND									
25	NOT FOUND									
26	NOT FOUND									
27	NOT FOUND									
28	NOT FOUND									
29	NOT FOUND									
30	NOT FOUND									
31	NOT FOUND									
32	NOT FOUND									
33	NOT FOUND									
34	NOT FOUND									
35	NOT FOUND									
36	NOT FOUND									
37	NOT FOUND									
38	NOT FOUND									
39	NOT FOUND									
40	91	608	15:19	3	0.696	A BB	855.	0.309 NG	0.11	
41	NOT FOUND									
42	NOT FOUND									
43	NOT FOUND									
44	NOT FOUND									
45	NOT FOUND									
46	NOT FOUND									
47	NOT FOUND									
48	NOT FOUND									
49	NOT FOUND									

PROCEDURE: FILTER/TIC DIAGNOSTIC REPORT 9/29/94 11:39:31

DATA FILE: CSLK929A

FILTER SCAN PARAMETERS

MAX. NUMBER TICS:	15
I1-TABLE ENTRIES:	539
SCAN TOLERANCE :	2
MIN. RIC HT (%):	10
FIRST SCAN :	1
LAST SCAN :	1600
TIC THRESHOLD :	600

METHOD LIBRARY & LISTS

TIC I.S. LIBRARY:	LIBRARYLS
NBS SEARCH PROC :	SERLIB
PEAK FINDER PROC:	VOME
TCA I.S. LL :	LS
FILE NAME LIST :	TCAREF2

TARGET COMPOUND ANALYSIS:

TARGETS (QUAN LIST)	IS PEAKS	TOTAL TARGET PEAKS
4	3	7

FILTER PROCESSING:

	<----------REJECT PEAKS---------->						
TOTAL PEAKS	< 1ST SCAN	> LAST SCAN	< MIN RIC HT	< SCAN TOL	> MAX # PEAKS	TOTAL REJECTS	TOTAL TICS
7	0	0	0	6	0	6	1

TIC PROCESSING:

NO.	SCAN#	PURITY	FIT	MW	COMPOUND NAME [BEFORE TIC THRESHOLD]
1	798	897	964	174	BENZENE, 1-BROMO-3-FLUORO-

Quantitation Report File: CBLK929A

Data: CBLK9[illegible] TI
[illegible]/29/94 [illegible]:00
Sample: M.[illegible]
Conds.: EPA METHOD 8240
Formula: SML Instrument: FINN Weight: 0.000
Submitted [illegible] PTL Analyst: UC Acct. No.:

AMOUNT=AREA * REF AMNT/(REF AREA * RESP FACT)
Resp. fac. from Library Entry

No	CAS	Name
1	[illegible]	CI01 BROMOCHLOROMETHANE **INT. STD. **
2	[illegible]	CI10 1,4-DIFLUOROBENZENE **INT. STD. **
3	[illegible]	CI20 CHLOROBENZENE-D5 **INT. STD. **
4	1073[illegible]	BENZENE, 1-BROMO-3-FLUORO-

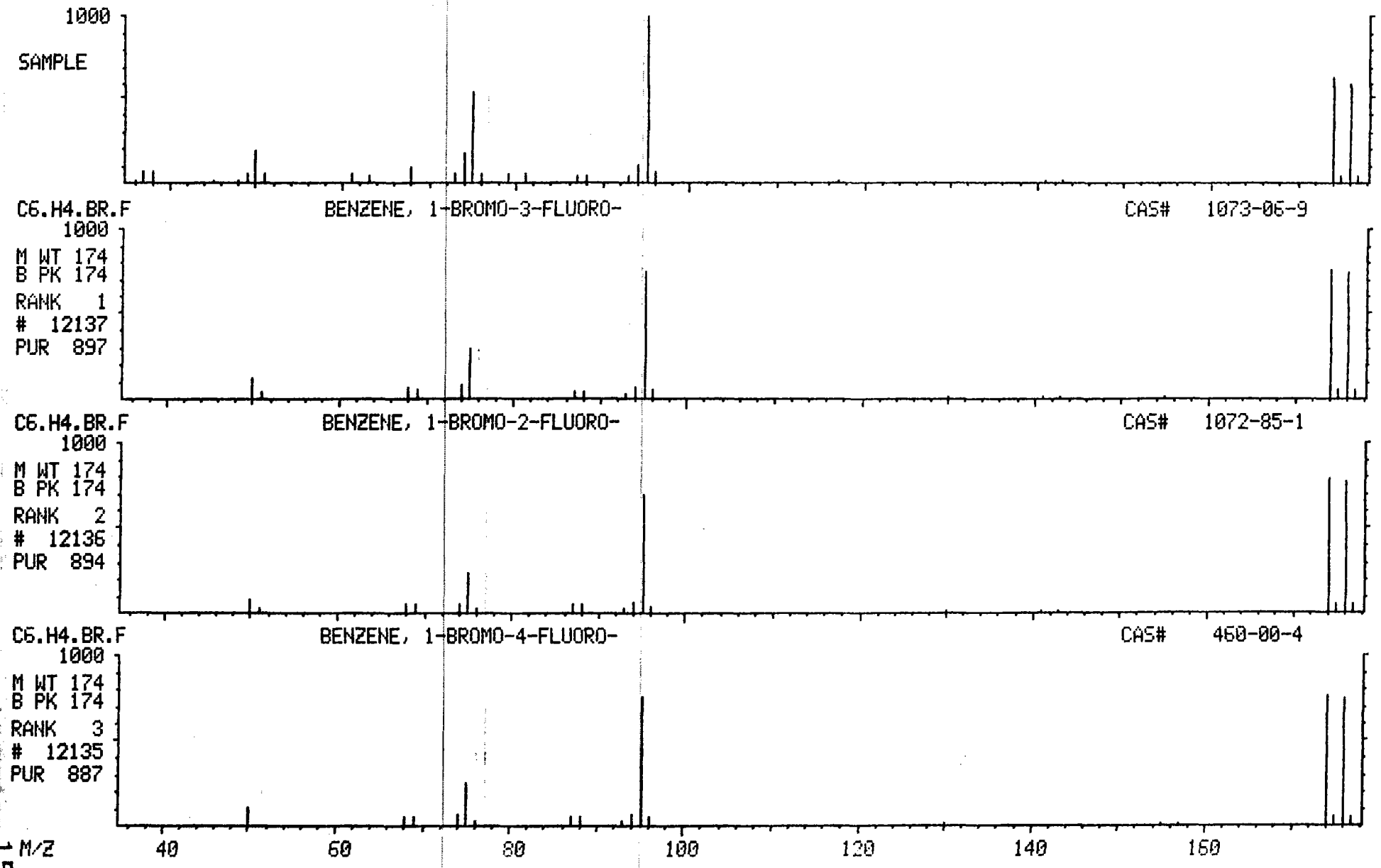
No	m/z	Base	Time	Ref	RRT	Meth	Area(Hght)	Amount		%Tot
1	TOT	[illegible]	7:32	0	ISINV	A BB	33933.	******	UG/L	00.00
2	TOT	[illegible]	9:48	0	ISINV	A BB	266301.	******	UG/L	00.00
3	TOT	[illegible]	21:59	0	ISINV	A BB	229777.	******	UG/L	00.00
4	TOT	[illegible]	20:06	3	0.914	A BB	256698.	55.858		*100.

No	Ret(L.	Ratio	RRT(L)	Ratio	Amnt	Amnt(L)	R.Fac	R.Fac(L)	Ratio
1	9:13	82	1.000						
2	19:21	[illegible].51	1.000						
3	23:54	92	1.000						
4					55.86	1.00	55.858	1.000	55.86

MID LIBRARY SEARCH (LIBRARYNB)
09/29/94 10:26:00 + 20:06
SAMPLE: M.BLK
CONDS.: EPA METHOD 8240
ENHANCED (S 15B 2N 0T)

DATA: CBLK929A # 798
CALI: CBLK929A # 3

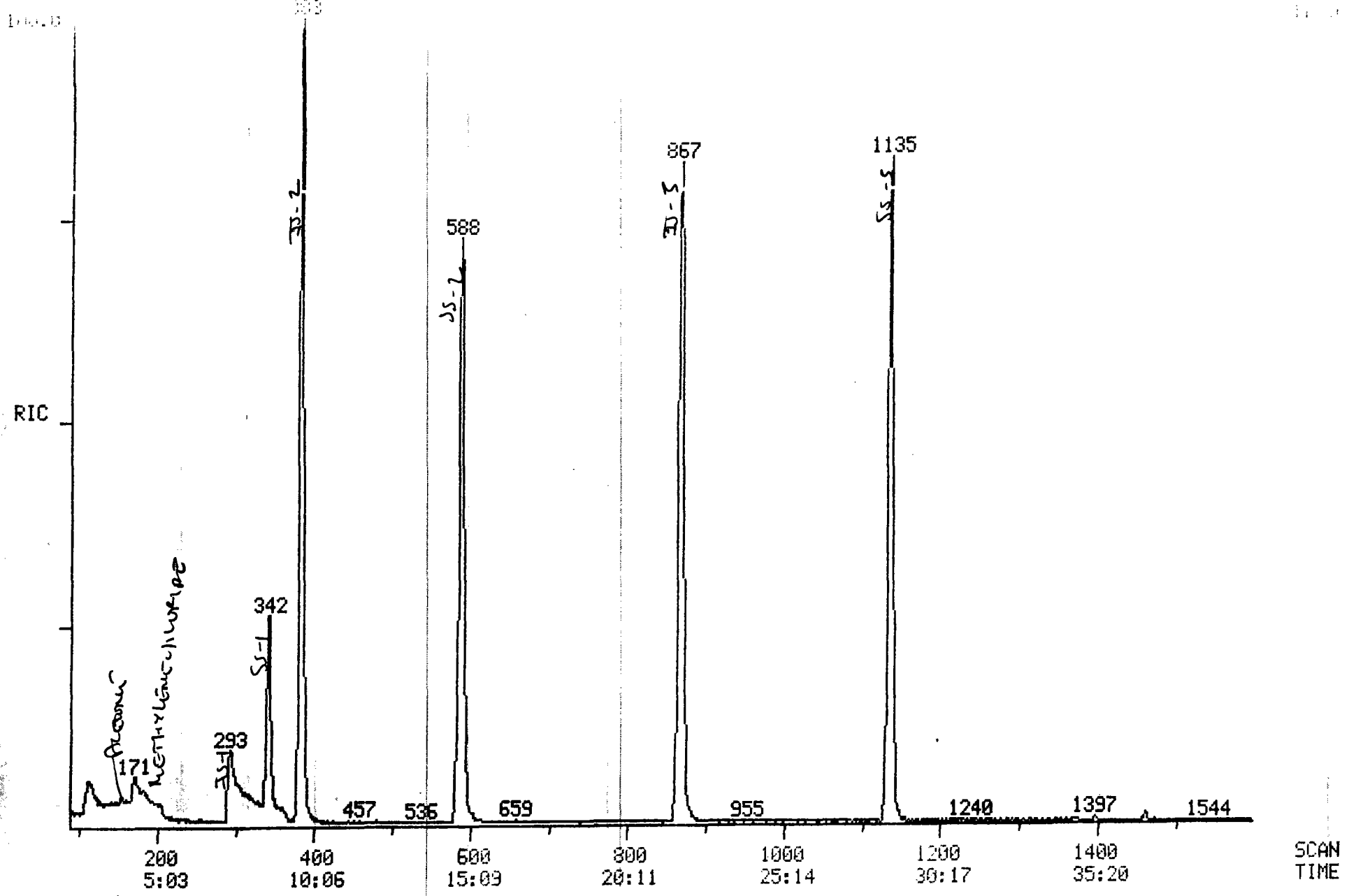
BASE M/Z: 95
RIC: 12192.



RIC
10/04/94 13:33:00
SAMPLE: M.BLK
CONDS.: EPA METHOD 8240
RANGE: G 1.1597 LABEL: N 0. 4.0 QUAN: N 0. 1.0 I 0 0.0E+00 0

DATA: CBLK1004 #1
CALI: CBLK1004 #3

SCANS 88 TO 1597



Quantitation Report File: CBLK1004

Data: CBLK1004.TI
09/04/94 13:33:00
Sample: M.BLK
Meths.: EPA METHOD 8240
Formula: Instrument: FINN Weight: 0.000
Submitted by: PTL Analyst: UC Acct. No.: -

AMOUNT=AREA * REF AMNT/(REF AREA * RESP FACT)
Resp. fac. from Library Entry

No	Name
1	CI01 BROMOCHLOROMETHANE **INT. STD. **
2	CI10 1,4-DIFLUOROBENZENE **INT. STD. **
3	CI20 CHLOROBENZENE-D5 **INT. STD. **
4	CS15 1,2-DICHLOROETHANE-D4 **S. STD. **
5	CS05 TOLUENE-D8 **S. STD. **
6	CS10 4-BROMOFLUOROBENZENE **S. STD. **
7	C010 CHLOROMETHANE **
8	C015 BROMOMETHANE
9	C020 VINYL CHLORIDE *
10	C025 CHLOROETHANE
11	C030 METHYLENE CHLORIDE
12	C251 ACROLIN
13	C035 ACETONE
14	C252 ACRYLONITRILE
15	C040 CARBON DISULFIDE
16	C045 1,1-DICHLOROETHENE *
17	C050 1,1-DICHLOROETHANE **
18	C055 TRANS-1,2-DICHLOROETHENE
19	C000 TRICHLOROFLUOROMETHANE
20	C060 CHLOROFORM *
21	C065 1,2-DICHLOROETHANE
22	C110 2-BUTANONE
23	C115 1,1,1-TRICHLOROETHANE
24	C120 CARBON TETRACHLORIDE
25	C125 VINYL ACETATE
26	C130 BROMO DICHLOROMETHANE
27	C140 1,2-DICHLOROPROPANE *
28	C145 TRANS-1,3 DICHLOROPROPENE
29	C150 TRICHLOROETHENE
30	C155 DIBROMOCHLOROMETHANE
31	C160 1,1,2-TRICHLOROETHANE
32	C165 BENZENE
33	C143 CIS-1,3-DICHLOROPROPENE
34	C175 2-CHLOROETHYL VINYL ETHER
35	C180 BROMOFORM **
36	C220 TETRACHLOROETHENE
37	C210 2-HEXANONE
38	C205 4-METHYL 2-PENTANONE
39	C225 1,1,2,2-TETRACHLOROETHANE **
40	C230 TOLUENE *
41	C235 CHLOROBENZENE **
42	C240 ETHYL BENZENE *
43	C245 STYRENE
44	C250 M+P-XYLENES
45	C253 1,3-DICHLOROBENZENE
46	C254 1,4-DICHLOROBENZENE
47	C255 1,2-DICHLOROBENZENE

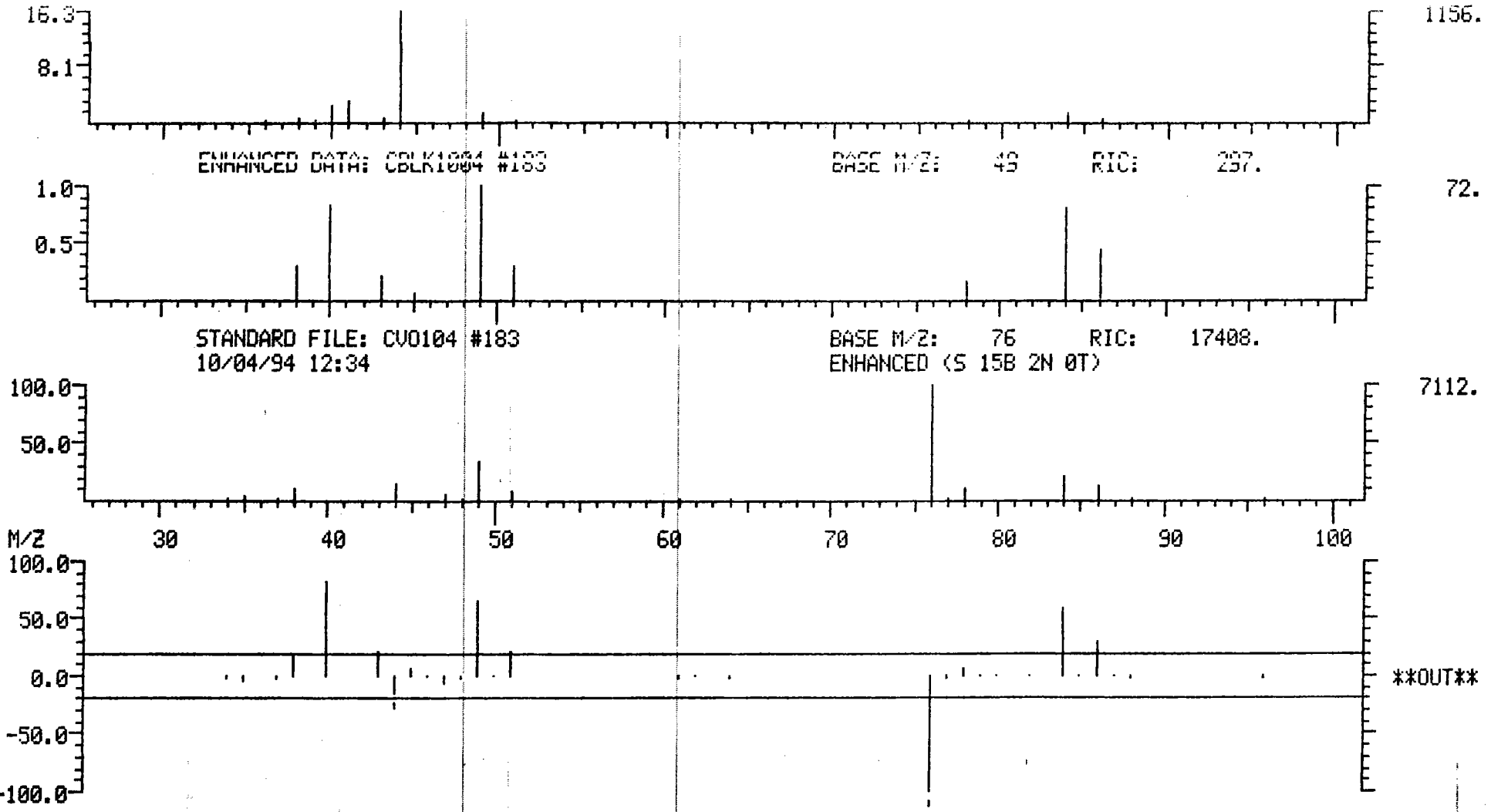
CDUK1004

No	Name
48	C250 O-XYLENE
49	BENZENE-D6 **S. STD. **

No	m/z	Scan	Time	Ref	RRT	Meth	Area(Hght)	Amount	%Tot
1	49	293	7:24	1	1.000	M XX	10934.	50.000 NG	16.48
2	114	382	9:38	2	1.000	A BB	102320.	50.000 NG	16.48
3	117	867	21:53	3	1.000	A BB	53294.	50.000 NG	16.48
4	65	341	8:36	1	1.164	A BB	17544.	44.054 NG	14.52
5	98	588	14:50	3	0.678	A BB	71077.	54.167 NG	17.85
6	95	1135	28:39	3	1.309	A BB	53377.	51.802 NG	17.07
7	NOT FOUND								
8	NOT FOUND								
9	NOT FOUND								
10	NOT FOUND								
11	49	182	4:36	1	0.621	A BB	480.	0.667 NG	0.22
12	NOT FOUND								
13	43	156	3:56	1	0.532	M XX	749.	2.560 NG	0.84
14	NOT FOUND								
15	NOT FOUND								
16	NOT FOUND								
17	NOT FOUND								
18	NOT FOUND								
19	NOT FOUND								
20	NOT FOUND								
21	NOT FOUND								
22	NOT FOUND								
23	NOT FOUND								
24	NOT FOUND								
25	NOT FOUND								
26	NOT FOUND								
27	NOT FOUND								
28	NOT FOUND								
29	NOT FOUND								
30	NOT FOUND								
31	NOT FOUND								
32	NOT FOUND								
33	NOT FOUND								
34	NOT FOUND								
35	NOT FOUND								
36	NOT FOUND								
37	NOT FOUND								
38	NOT FOUND								
39	NOT FOUND								
40	91	601	15:10	3	0.693	A BB	587.	0.881 NG	0.13
41	NOT FOUND								
42	NOT FOUND								
43	NOT FOUND								
44	NOT FOUND								
45	NOT FOUND								
46	NOT FOUND								
47	NOT FOUND								
48	NOT FOUND								
49	NOT FOUND								

DATA FILE: CBLK1004 #183 STANDARD FILE: CV0104 #183
TARGET COMPOUND COMPARISON CALI: CBLK1004 #3
COMPOUND: C030 METHYLENE CHLORIDE

RAW DATA: CBLK1004 #183 BASE M/Z: 49 RIC: 1972.
10/04/94 13:33

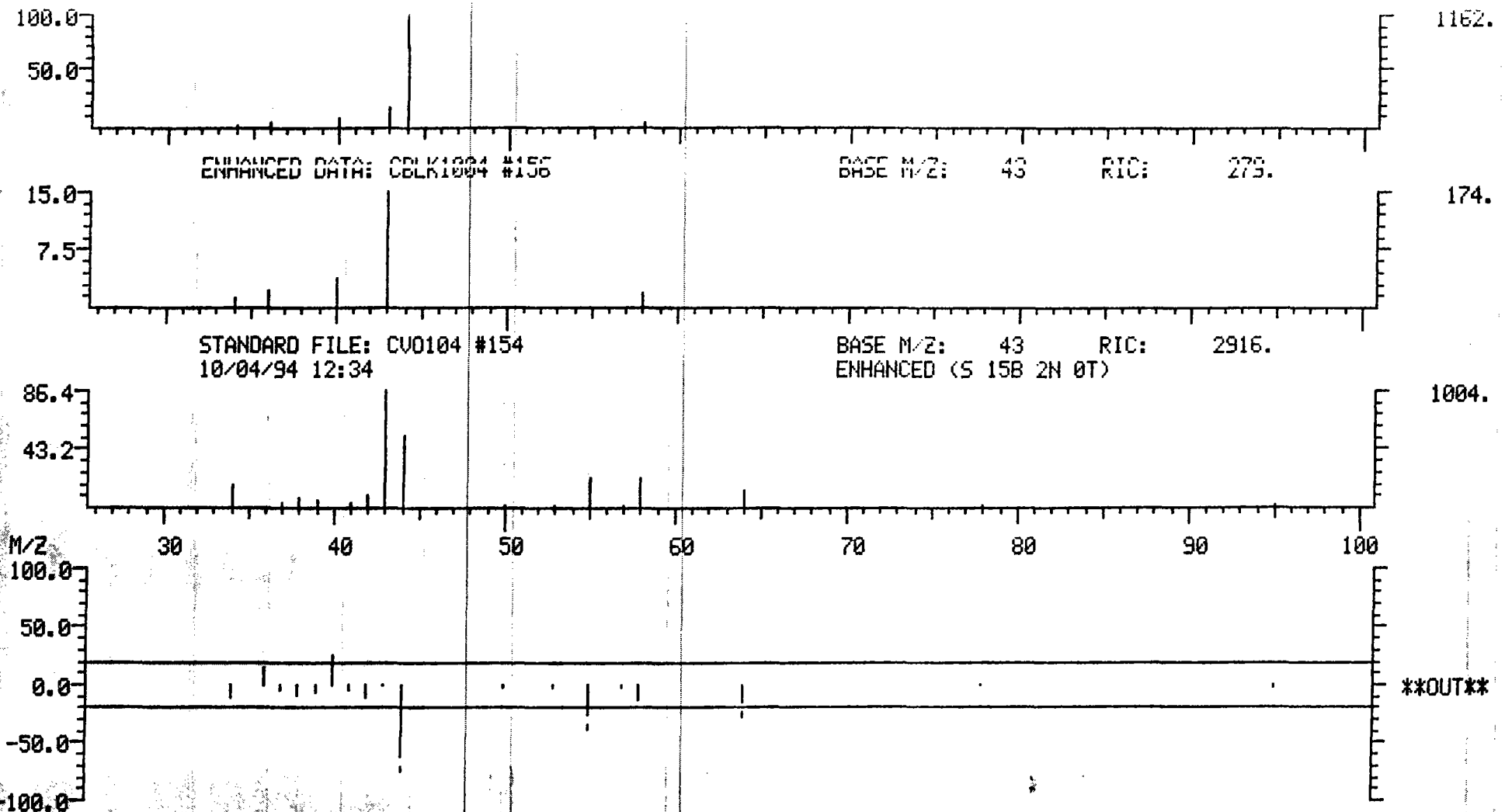


DATA FILE: CBLK1004 #156
TARGET COMPOUND COMPARISON
COMPOUND: C035 ACETONE

STANDARD FILE: CV0104 #154
CALI: CBLK1004 #3

RAW DATA: CBLK1004 #156
10/04/94 13:33

BASE M/Z: 44 RIC: 1574.



PROCEDURE: ...TER/TIC DIAGNOSTIC REPORT 10/05/94 8:28:28

DATA FILE: ...A1004

FILTER SCA... PARAMETERS

MAX. NUMBE... PKS:	15
11-TABLE E... ES:	539
SCAN TOLE... :	2
MIN. RIC ... T:	10
FIRST SCAN :	1
LAST SCAN :	1600
TIC THRESH... :	600

METHOD LIBRARY & LISTS

TIC I.S. LIBRARY:	LIBRARYLS
NBS SEARCH PROC :	SERLIB
PEAK FINDER PROC:	VOME
TCA I.S. LL :	LS
FILE NAME LIST :	TCAREF2

TARGET COM... D ANALYSIS:

TARGETS (QUAN LIST)	IS PEAKS	TOTAL TARGET PEAKS
6	3	9

FILTER PRO... ING:

TOTAL PEAKS	...	REJECT PEAKS: > LAST SCAN	< MIN RIC HT	< SCAN TOL	> MAX # PEAKS	TOTAL REJECTS	TOTAL TICS
8		0	1	6	0	7	1

...C PROCES...

NO.	SCAN...	...URITY	FIT	MW	COMPOUND NAME [BEFORE TIC THRESHOLD]
1	...	410	427	85	ACETIC ACID, CYANO-

Quantitat[illegible]eport File: CBLK1004

Data: CBLK[illegible].TI
[illegible]/04/94 [illegible].00
Sample: M.[illegible]
Conds.: EP[illegible]THOD 8240
Formula: Instrument: FINN Weight: 0.000
Submitted [illegible]TL Analyst: UC Acct. No.: -

AMOUNT=AREA [illegible]/EF AMNT/(REF AREA * RESP FACT)
Resp. fac. [illegible]m Library Entry

No	CAS	Name
1	[illegible]-0	CI01 BROMOCHLOROMETHANE **INT. STD. **
2	[illegible]-0	CI10 1,4-DIFLUOROBENZENE **INT. STD. **
3	[illegible]-0	CI20 CHLOROBENZENE-D5 **INT. STD. **
4	[illegible]-0	UNKNOWN

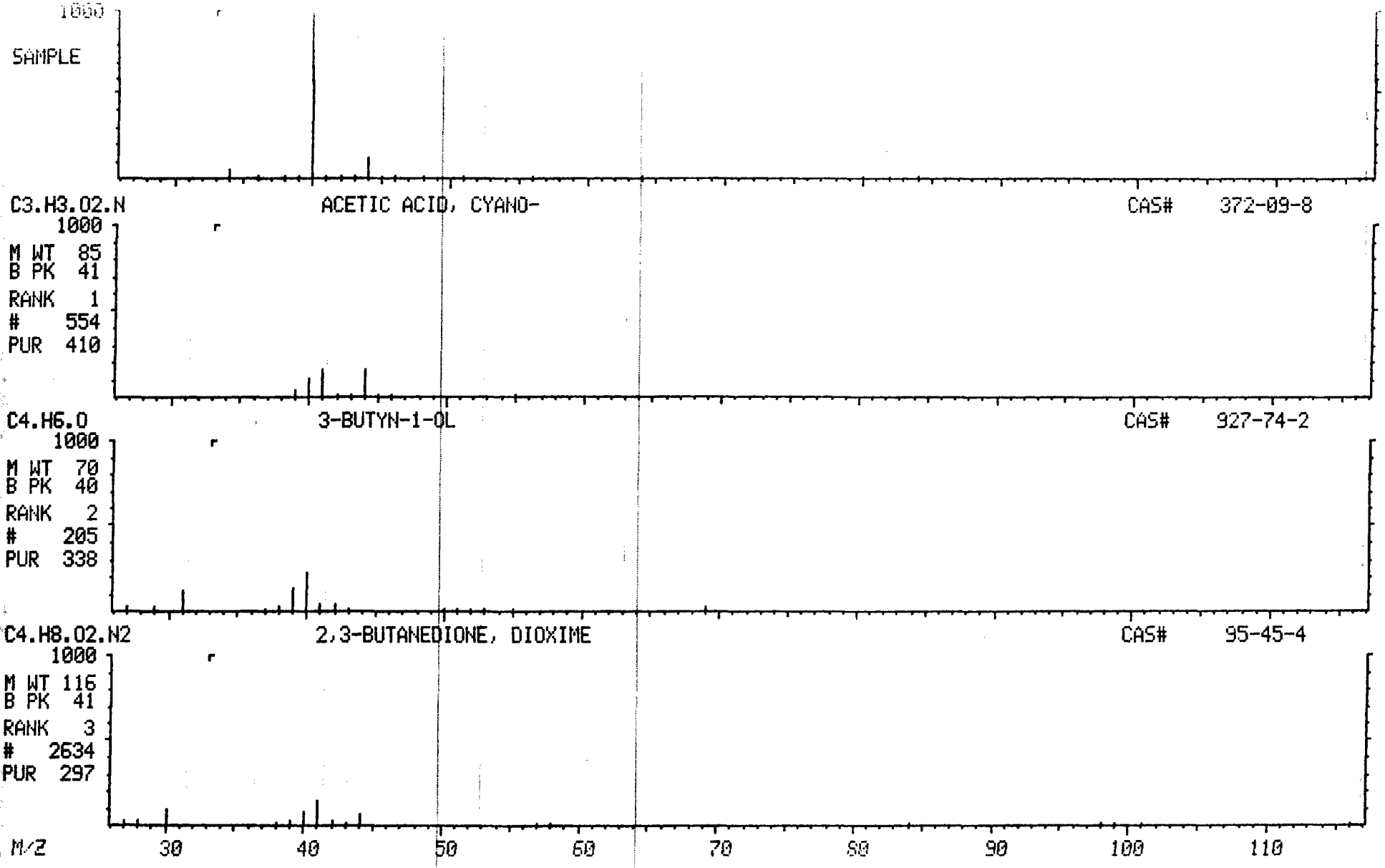
No	m/z	[illegible]	Time	Ref	RRT	Meth	Area(Hght)	Amount		%Tot
1	TOT	[illegible]	7:22	0	ISINV	A BB	26158.	******	UG/L	00.00
2	TOT	[illegible]	9:40	0	ISINV	A BB	217284.	******	UG/L	00.00
3	TOT	[illegible]	21:53	0	ISINV	A BB	183485.	******	UG/L	00.00
4	TOT	[illegible]	2:01	1	0.274	A BB	641968.	1227.100		*100.

No	Ret(L)	[illegible]atio	RRT(L)	Ratio	Amnt	Amnt(L)	R.Fac	R.Fac(L)	Ratio
1	9:13	[illegible]80	1.000						
2	19:21	[illegible]50	1.000						
3	23:54	[illegible]92	1.000						
4					1227.10	1.001227.100		1.0001227.10	

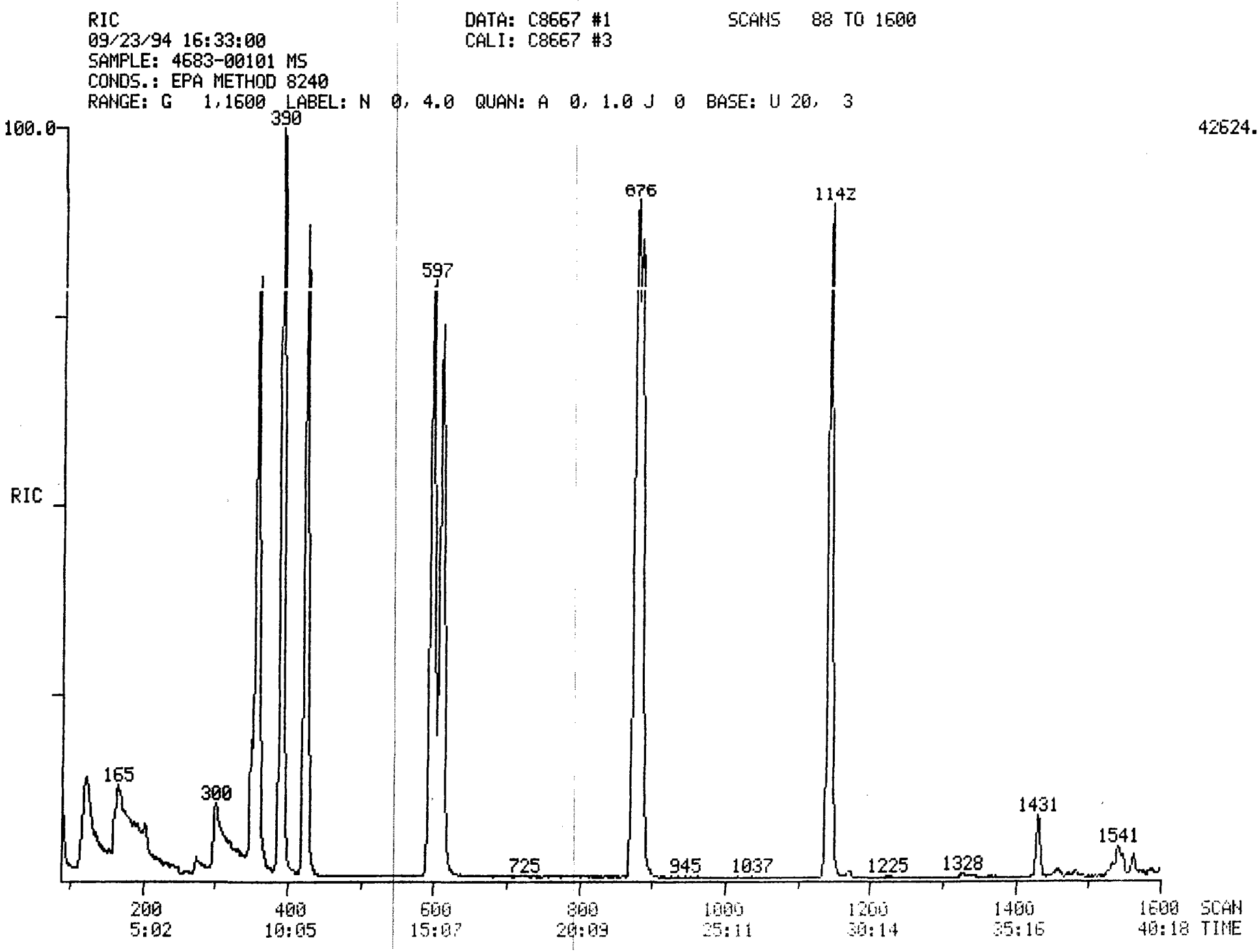
MID LIBRARY SEARCH (LIBRARYNB)
10/04/94 13:33:00 + 2:01
SAMPLE: M.BLK
CONDS.: EPA METHOD 8240
ENHANCED (S 15B 2N 0T)

DATA: CBLK1004 # 80
CALI: CBLK1004 # 3

BASE M/Z: 40
RIC: 241664.



167


RIC
09/23/94 16:33:00
SAMPLE: 4683-00101 MS
CONDS.: EPA METHOD 8240
RANGE: G 1,1600 LABEL: N 0, 4.0 QUAN: A 0, 1.0 J 0 BASE: U 20, 3
DATA: C8667 #1
CALI: C8667 #3
SCANS 88 TO 1600
100.0
42624.
RIC
390
676
1142
597
165
300
725
945
1037
1225
1328
1431
1541
200
5:02
400
10:05
600
15:07
800
20:09
1000
25:11
1200
30:14
1400
35:16
1600
40:18
SCAN
TIME

Quantitation Report File: C8667

Data: C8667.TI
?/23/94 16:33:00
Sample: 4683-00101 MS
Conds.: EPA METHOD 8240
Formula: 2.5G/5ML Instrument: FINN Weight: 0.000
Submitted by: USARMY Analyst: UC Acct. No.: 4683-001

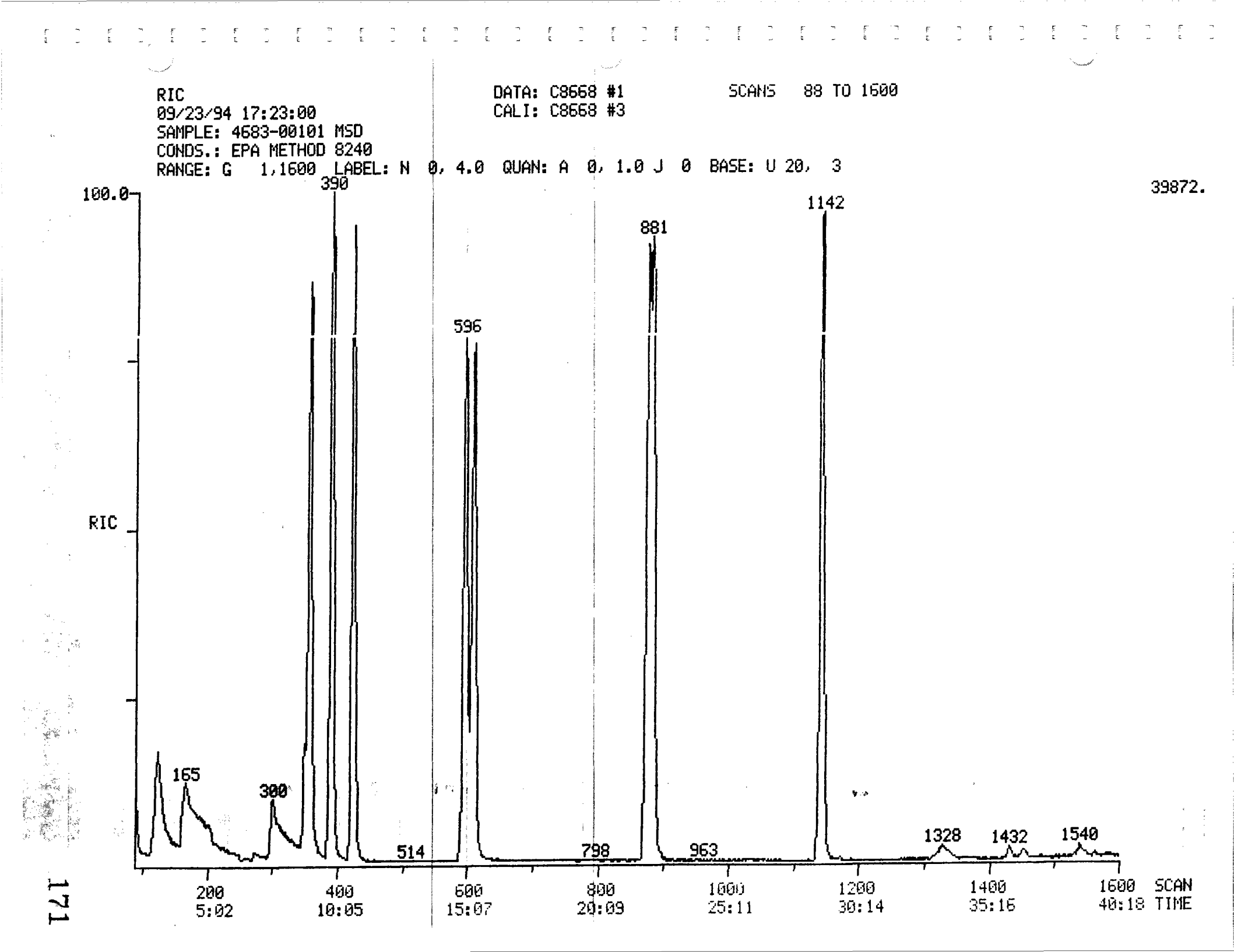
AMOUNT=AREA * REF AMNT/(REF AREA * RESP FACT)
Resp. fac. from Library Entry

No	Name
1	CI01 BROMOCHLOROMETHANE **INT. STD. **
2	CI10 1,4-DIFLUOROBENZENE **INT. STD. **
3	CI20 CHLOROBENZENE-D5 **INT. STD. **
4	CS15 1,2-DICHLOROETHANE-D4 **S. STD. **
5	CS05 TOLUENE-D8 **S. STD. **
6	CS10 4-BROMOFLUOROBENZENE **S. STD. **
7	C010 CHLOROMETHANE **
8	C015 BROMOMETHANE
9	C020 VINYL CHLORIDE *
10	C025 CHLOROETHANE
11	C030 METHYLENE CHLORIDE
12	C251 ACROLIN
13	C035 ACETONE
14	C252 ACRYLONITRILE
15	C040 CARBON DISULFIDE
16	C045 1,1-DICHLOROETHENE *
17	C050 1,1-DICHLOROETHANE **
18	C055 TRANS-1,2-DICHLOROETHENE
19	C000 TRICHLOROFLUOROMETHANE
20	C060 CHLOROFORM *
21	C065 1,2-DICHLOROETHANE
22	C110 2-BUTANONE
23	C115 1,1,1-TRICHLOROETHANE
24	C120 CARBON TETRACHLORIDE
25	C125 VINYL ACETATE
26	C130 BROMO DICHLOROMETHANE
27	C140 1,2-DICHLOROPROPANE *
28	C145 TRANS-1,3 DICHLOROPROPENE
29	C150 TRICHLOROETHENE
30	C155 DIBROMOCHLOROMETHANE
31	C160 1,1,2-TRICHLOROETHANE
32	C165 BENZENE
33	C143 CIS-1,3-DICHLOROPROPENE
34	C175 2-CHLOROETHYL VINYL ETHER
35	C180 BROMOFORM **
36	C220 TETRACHLOROETHENE
37	C210 2-HEXANONE
38	C205 4-METHYL 2-PENTANONE
39	C225 1,1,2,2-TETRACHLOROETHANE **
40	C230 TOLUENE *
41	C235 CHLOROBENZENE **
42	C240 ETHYL BENZENE *
43	C245 STYRENE
44	C250 M+P-XYLENES
45	C253 1,3-DICHLOROBENZENE
46	C254 1,4-DICHLOROBENZENE
47	C255 1,2-DICHLOROBENZENE

My 17

No	Name
48	C250 O-XYLENE
49	BENZENE-D6 **S. STD. **

No	m/z	Scan	Time	Ref	RRT	Meth	Area(Hght)	Amount	%Tot
1	49	300	7:33	1	1.000	A BB	8847.	50.000 NG	10.12
2	114	390	9:49	2	1.000	A BB	101972.	50.000 NG	10.12
3	117	875	22:02	3	1.000	A BB	74708.	50.000 NG	10.12
4	65	349	8:47	1	1.163	M XX	15979.	35.725 NG	7.23
5	98	597	15:02	3	0.682	A BB	93124.	46.564 NG	9.42
6	95	1142	28:46	3	1.305	A BB	59844.	38.727 NG	7.84
7	NOT FOUND								
8	NOT FOUND								
9	NOT FOUND								
10	NOT FOUND								
11	49	189	4:46	1	0.630	A BV	1220.	1.903 NG	0.39
12	NOT FOUND								
13	NOT FOUND								
14	NOT FOUND								
15	NOT FOUND								
16	96	166	4:11	1	0.553	A BB	5724.	44.136 NG	8.93
17	NOT FOUND								
18	NOT FOUND								
19	NOT FOUND								
20	NOT FOUND								
21	NOT FOUND								
22	NOT FOUND								
23	NOT FOUND								
24	NOT FOUND								
25	NOT FOUND								
26	NOT FOUND								
27	NOT FOUND								
28	NOT FOUND								
29	95	424	10:41	2	1.087	A BB	39899.	45.065 NG	9.12
30	NOT FOUND								
31	NOT FOUND								
32	78	357	9:00	2	0.915	A BB	86662.	44.339 NG	8.97
33	NOT FOUND								
34	NOT FOUND								
35	NOT FOUND								
36	NOT FOUND								
37	NOT FOUND								
38	NOT FOUND								
39	NOT FOUND								
40	91	610	15:22	3	0.697	A BB	90072.	41.129 NG	8.32
41	112	882	22:13	3	1.008	A BB	65297.	46.658 NG	9.44
42	NOT FOUND								
43	NOT FOUND								
44	NOT FOUND								
45	NOT FOUND								
46	NOT FOUND								
47	NOT FOUND								
48	NOT FOUND								
49	NOT FOUND								


RIC
09/23/94 17:23:00
SAMPLE: 4683-00101 MSD
CONDS.: EPA METHOD 8240
RANGE: G 1,1600 LABEL: N 0, 4.0 QUAN: A 0, 1.0 J 0 BASE: U 20, 3
DATA: C8668 #1
CALI: C8668 #3
SCANS 88 TO 1600
39872.
100.0
RIC
390
1142
881
596
165
300
514
798
963
1328
1432
1540
200
5:02
400
10:05
600
15:07
800
20:09
1000
25:11
1200
30:14
1400
35:16
1600
40:18
SCAN
TIME
171

Quantitation Report File: C8668

Data: C8668.TI
9/23/94 17:23:00
Sample: 4683-00101 MSD
Conds.: EPA METHOD 8240
Formula: 2.5G/5ML Instrument: FINN Weight: 0.000
Submitted by: USARMY Analyst: UC Acct. No.: 4683-001

AMOUNT=AREA * REF AMNT/(REF AREA * RESP FACT)
Resp. fac. from Library Entry

No	Name
1	CI01 BROMOCHLOROMETHANE **INT. STD. **
2	CI10 1,4-DIFLUOROBENZENE **INT. STD. **
3	CI20 CHLOROBENZENE-D5 **INT. STD. **
4	CS15 1,2-DICHLOROETHANE-D4 **S. STD. **
5	CS05 TOLUENE-D8 **S. STD. **
6	CS10 4-BROMOFLUOROBENZENE **S. STD. **
7	C010 CHLOROMETHANE **
8	C015 BROMOMETHANE
9	C020 VINYL CHLORIDE *
10	C025 CHLOROETHANE
11	C030 METHYLENE CHLORIDE
12	C251 ACROLIN
13	C035 ACETONE
14	C252 ACRYLONITRILE
15	C040 CARBON DISULFIDE
16	C045 1,1-DICHLOROETHENE *
17	C050 1,1-DICHLOROETHANE **
18	C055 TRANS-1,2-DICHLOROETHENE
19	C000 TRICHLOROFLUOROMETHANE
20	C060 CHLOROFORM *
21	C065 1,2-DICHLOROETHANE
22	C110 2-BUTANONE
23	C115 1,1,1-TRICHLOROETHANE
24	C120 CARBON TETRACHLORIDE
25	C125 VINYL ACETATE
26	C130 BROMO DICHLOROMETHANE
27	C140 1,2-DICHLOROPROPANE *
28	C145 TRANS-1,3 DICHLOROPROPENE
29	C150 TRICHLOROETHENE
30	C155 DIBROMOCHLOROMETHANE
31	C160 1,1,2-TRICHLOROETHANE
32	C165 BENZENE
33	C143 CIS-1,3-DICHLOROPROPENE
34	C175 2-CHLOROETHYL VINYL ETHER
35	C180 BROMOFORM **
36	C220 TETRACHLOROETHENE
37	C210 2-HEXANONE
38	C205 4-METHYL 2-PENTANONE
39	C225 1,1,2,2-TETRACHLOROETHANE **
40	C230 TOLUENE *
41	C235 CHLOROBENZENE **
42	C240 ETHYL BENZENE *
43	C245 STYRENE
44	C250 M+P-XYLENES
45	C253 1,3-DICHLOROBENZENE
46	C254 1,4-DICHLOROBENZENE
47	C255 1,2-DICHLOROBENZENE

MSD 61

No	Name
48	C250 O-XYLENE
49	BENZENE-D6 **S. STD. **

No	m/z	Scan	Time	Ref	RRT	Meth	Area(Hght)	Amount	%Tot
1	49	300	7:33	1	1.000	A BB	8993.	50.000 NG	9.93
2	114	390	9:49	2	1.000	A BB	96476.	50.000 NG	9.93
3	117	875	22:02	3	1.000	A BB	72942.	50.000 NG	9.93
4	65	349	8:47	1	1.163	M XX	16470.	36.225 NG	7.19
5	98	596	15:01	3	0.681	A BB	86242.	44.167 NG	8.77
6	95	1142	28:46	3	1.305	A BB	59490.	39.431 NG	7.83
7	NOT FOUND								
8	NOT FOUND								
9	NOT FOUND								
10	NOT FOUND								
11	NOT FOUND								
12	NOT FOUND								
13	NOT FOUND								
14	NOT FOUND								
15	NOT FOUND								
16	96	165	4:09	1	0.550	A BB	5812.	44.087 NG	8.75
17	NOT FOUND								
18	NOT FOUND								
19	NOT FOUND								
20	NOT FOUND								
21	NOT FOUND								
22	NOT FOUND								
23	NOT FOUND								
24	NOT FOUND								
25	NOT FOUND								
26	NOT FOUND								
27	NOT FOUND								
28	NOT FOUND								
29	95	423	10:39	2	1.085	A BB	40842.	48.758 NG	9.68
30	NOT FOUND								
31	NOT FOUND								
32	78	356	8:58	2	0.913	A BB	87207.	47.160 NG	9.36
33	NOT FOUND								
34	NOT FOUND								
35	NOT FOUND								
36	NOT FOUND								
37	NOT FOUND								
38	NOT FOUND								
39	NOT FOUND								
40	91	609	15:20	3	0.696	A BB	92584.	43.300 NG	8.60
41	112	882	22:13	3	1.008	A BB	69143.	50.603 NG	10.05
42	NOT FOUND								
43	NOT FOUND								
44	NOT FOUND								
45	NOT FOUND								
46	NOT FOUND								
47	NOT FOUND								
48	NOT FOUND								
49	NOT FOUND								

After the repairing of maintainance

	D/T ANALYZED	DATA FILE	JOB NUMBER	PTL ID #	CLIENT	DILUTION	NON	SURROGATE Method 8240	COMMENTS	INITS
	08/14/94	BFB914	Paired @ 12:00.							
9		✓CV914SP		50 ppb	NOA std"	SPL - Method 8240 Std.			#77 no surr.d	
3		✓CV91420		20 ppb	NOA std				#34 Missing	
3		✓CV91410		10 ppb	NOA std		34M		#34M	
3		✓CV914	100A	100 ppb	NOA std				#34M	
4		CV914	200 ✓	200 ppb	NOA std		#34			
3		✓CV914	50	50 ppb	NOA std				Good	pc
5		CV914	100B ✓	100 ppb						
6		CV914	50A ✓	50 ppb						
7		CV914	20A ✓	20 ppb						

	D/T ANALYZED	DATA FILE	JOB NUMBER	PTL ID #	CLIENT	DILUTION	NON	SURROGATE Method 8240	COMMENTS	INITS
	9/23/94	BFB923	Paired @	IC:51		10:55			142.151	
3		CV0923	50	CAB	Vol Std				ok	
4		CBLK923	M. Blank						clean ok	
5		C8663	4714-	00101	Zenith	0.5/125 10ml/5ml:	x125 dil.		ok #4	
6		C8664	4682-	00101	US Army	5ml			ok	
7		C8665	4683-	00101	↓	↓			ok	
8		C8666	4747-	01	Environmental	2.5g/5ml			ok	
9		C8667	4683-	MS 00101	US Army	↓		940923-SY	ok	
10		C8668		MSD 00101	↓	↓			ok	
1		C8669	4601-	RR 00301	Delcora	5S/10ml ÷2 10ml/5ml = 10	x10 dil		ok	
2		C8670	-	RR 00401	↓	↓			ok	

	D/T ANALYZED	DATA FILE	JOB NUMBER	PTL ID	CLIENT	DILUTION	IS NON	SURROGATE Meas'd 8240	COMMENTS	INITS
	9/29/94	BFB929		Laird @ 8:45					1H8, 150	VL
3		CV0929		50 ppb	VOA std				OK	VL
4		CB1-939A		M: Blank					OK not loaded	VL
		C8686	4632-	TCLP Blk		1:5		Same as Lab Hydro did not run		VL
5		C8687	4632-	TCLP 01	Briggs				OK	VL
6		C8688	-	TCLP 02					OK	VL
7		C8689	4508-	TCLP 01	NAVCAD				OK	VL
8		C8690	-	TCLP 02					OK	VL
9		C8691	-	TCLP 03					OK	VL
10		C8692	-	TCLP 04					OK	VL
1		C8693	4339-	TCLP 01	Eastern					

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D/T ANALYZED	DATA FILE	JOB NUMBER	PTL ID	CLIENT	DILUTION	IS KOH	SURROGATE	COMMENTS	UNITS
9/24/97					5g/10ml = 2				
2	C8694	4760-	01	US Army	100ml/5ml	50		OK	ug
3	C8695	-	02	↓	↓			OK	ug
					5g/10ml = 2				
4	C8696	4760	02		200ml/5ml	35		OK	ug
5	C8697	-	01					OK	ug
6	C8698	4755-	02	CW	5ml			OK	ug
			MST44						
7	C8699	4632-	01	Briggs	1:5			OK	ug
9/20/97 3	C8700	4820		RSW88	10ml	Aq	Sample		

	D/T ANALYZED	DATA FILE	JOB NUMBER	PTL ID	CLIENT	DILUTION	IS NON	SURROGATE	COMMENTS	INITS
		BFB1004B		Paired @ 11:42					146,145	
3		CV01004		BPB	VOA std					
4		CBLK1004		Blank						
5		C8711	4760-	00103	US Army	5 ml				
6		C8712	4899-	00101	↓	5/5 ml	1:5			
7		C8713	4652-	13-RR 00101	Azmiler	Back 100 ml/100 ml 100 ml/5 ml x 10,000 dil				
8		C8714	—		↓	front 5 ml/5 ml				
9		C8715	4871	00102	CW	5g/10 ml F 100 ml/5 ml F 100 x dil				
10		C8716	4842	00103	Burlington	?.5g/5 ml	x2 dil			
1		C8717		MS	↓	↓				
2		C8718		MSD	↓	↓				

	D/T ANALYZED	DATA FILE	JOB NUMBER	PTL ID	CLIENT	DILUTION	IS NON	SURROGATE	COMMENTS	UNITS
				1651-1						
3	6/04/90	C8719	4899-	00101	US Army	2.5 5/5ml			"	"
4		C8720	-	00102	"	5ml			"	"
5		C8721	J.H2O-			5ml			"	"

US ARMY FT. MONMOUTH ENVIRONMENTAL LABORATORY
NJDEPE # 13461

REPORT OF ANALYSIS

Client: U.S. Army
DPW, SELFM-PW-EV
Bldg. 173
Ft. Monmouth, NJ 07703

Project: Total Petroleum Hydrocarbons
98-0001
Bldg.290

Project # 3437
Date Rec. 03/26/98
Date Compl. 03/27/98
Released by:

Daniel K. Wright
Laboratory Director

Table of Contents

Method Summary

NJDEP Method OQA-QAM-025-10/97

Gas Chromatographic Determination of Total Petroleum Hydrocarbons in Soil

Fifteen grams (15g)(wet weight) of a soil sample is added to a 125 mL acid cleaned, solvent rinsed, capped Erlenmeyer flask. 15g anhydrous sodium sulfate is added to dry sample. Surrogate standard spiking solution is then added to the flask.

Twenty five milliliters(25mL) Methylene Chloride is added to the flask and it is secured on a gyrotory shaker table. The agitation rate is set to 400rpm and the sample is shaken for 30 minutes. The flask is the removed from the table and the particulate matter is allowed to settle. The extract is transferred to a Teflon capped vial. A second 25mL of Methylene Chloride is added to the flask and shaken for an additional 30 minutes. The flask is again removed and allowed to settle. The extracts are combined in the vial then transferred to a 1mL autosampler vial.

The extract is then injected directly into a GC-FID for analysis. The sample is analyzed for petroleum hydrocarbons covering a range of C8-C42 including pristane and phytane. Total Petroleum Hydrocarbon concentration is determined by integrating between 5 minutes and 22 minutes. The baseline is established by starting the integration after the end of the solvent peak and stopping after the last peak.

The final concentration of Total Petroleum Hydrocarbons is calculated using percent solid, sample weight and concentration.

PHC Conformance/Non-conformance Summary Report

	No	Yes
1. Method Detection Limits provided.	__	✓
2. Method Blank Contamination - If yes, list the sample and the corresponding concentrations in each blank.	✓	__
3. Matrix Spike Results Summary Meet Criteria. (If not met, list the sample and corresponding recovery which falls outside the acceptable range).	__	✓
4. Duplicate Results Summary Meet Criteria. (If not met, list the sample and corresponding recovery which falls outside the acceptable range).	__	✓
5. IR Spectra submitted for standards, blanks, & samples	__	NA __
6. Chromatograms submitted for standards, blanks, and samples if GC fingerprinting was conducted.	__	✓
7. Analysis holding time met. (If not met, list number of days exceeded for each sample)	__	✓

Additional Comments: _____

Laboratory Authentication Statement

I certify under penalty of law, where applicable, that this laboratory meets the Laboratory Performance Standards and Quality Control requirements specified in N.J.A.C. 7:18 and 40 CFR Part 136 for Water and Wastewater Analyses and SW 846 for Solid Waste Analysis. I have personally examined the information contained in this report, and to the best of my knowledge, I believe that the submitted information is true, accurate, complete, and meets the above referenced standards where applicable. I am aware that there are significant penalties for purposefully submitting falsified information, including the possibility of a fine and imprisonment.

Daniel K. Wright
Laboratory Manager

Fort Monmouth Environmental Testing Laboratory

Bldg. 173, SELFM-PW-EV, Fort Monmouth, NJ 07703
Tel (732)532-4359 Fax (732)532-3484 EMail:appleby@doim6.monmouth.army.mil
NJDEP Certification #13461

Chain of Custody Record

Customer: [illegible]
Phone #: [illegible]
( )DERA (X)OMA ( )Other:_____

Project No: GST/98-0001
Location: Bldg-290

Samplers Name / Company :

Lab Sample I.D.	Sample Location	Date	Time	Sample Type	# bottles	VOA + 15	TPH							Comments: Remarks / Preservation Method
						Analysis Parameters								
3437.01	TRIP BLANK	3-26-98	~~1348~~	AQ	2	X								OVA #2-AS/1903 HCL
02	290-B-1		1348	SOIL	1		X							9 PPM) 24-30" (ACC)
03	290-B-2		1433	SOIL	1		X							100 PPM) 22-26" (MC)
04	290-B-3		1507	SOIL	1		X							90 PPM)
(MC) ~~05~~	~~290-B-4~~			~~SOIL~~	~~1~~		~~X~~							
05	290-B-2-W		1550	AQ	2	X								
06	290-B-4		1625	SOIL	1		X							

Relinquished by (signature):	Date/Time:	Received by (signature):	Relinquished by (signature):	Date/Time:	Received by (signature):
[signature]	3-26-98 1640	[signature]			
Relinquished by (signature):	Date/Time:	Received by (signature):	Relinquished by (signature):	Date/Time:	Received by (signature):
		[signature]			

Report Type: (_)Full, (_)Reduced, (_)Standard, (_)Screen / non-certified
Turnaround time: (_)Standard 4 wks, (_)Rush ____ Days, (_)ASAP Verbal ____ Hrs.

Remarks:

print legibly

~me1a23.xls2/20/98

Report of Analysis
U.S. Army, Fort Monmouth Environmental Laboratory
NJDEP Certification # 13461

Client :	U.S. Army	**Lab. ID # :**	3437
	DPW. SELFM-PW-EV	**Date Rec'd:**	26-Mar-98
	Bldg. 173	**Analysis Start:**	27-Mar-98
	Ft. Monmouth, NJ 07703	**Analysis Complete:**	27-Mar-98
Analysis:	OQA-QAM-025	**UST Reg. #:**	
Matrix:	Soil	**Closure #:**	
Analyst:	D.DEINHARDT	**DICAR #:**	
Ext. Meth:	Shake	**Location #:**	BLDG. 290

Sample	Field ID	Dilution Factor	Weight (g)	% Solid	MDL (mg/kg)	TPHC Result (mg/kg)
3437.02	290-B-1	1.00	14.98	92.81	169	317.67
3437.03	290-B-2	1.00	15.11	84.28	185	ND
3437.04	290-B-3	1.00	15.22	78.81	196	224.45
3437.06	290-B-4	1.00	16.04	74.89	196	ND
METHOD BLANK	27-Mar-98	1.00	15.00	100.00	157	ND

ND = Not Detected

MDL = Method Detection Limit

Daniel K. Wright
Laboratory Director

Response Factor Report FID/TCD

Method : C:\HPCHEM\1\METHODS\TPH27.M (Chemstation Integrator)
Title : TPHC Calibration 06/05/97 21 peaks
Last Update : Thu Mar 19 07:39:01 1998

Calibration Files
200 =T04649.D 100 =T04654.D 50 =T04651.D
10 =T04652.D 5 =T04653.D

		Compound	200	100	50	10	5	Avg		%RSD
1)	tC	C8	1.801	2.041	1.835	1.725	1.727	1.826	E4	7.10
2)	tC	C10	1.933	2.259	1.974	1.798	1.804	1.953	E4	9.60
3)	TC	C12	2.116	2.476	2.169	1.953	1.968	2.136	E4	9.89
4)	tC	C14	2.185	2.559	2.270	2.025	2.004	2.209	E4	10.19
5)	tC	C16	2.233	2.625	2.327	2.097	2.075	2.271	E4	9.81
6)	tC	C18	2.504	2.926	2.710	2.444	2.375	2.592	E4	8.67
7)	tC	C20	2.415	2.851	2.515	2.261	2.228	2.454	E4	10.22
8)	tC	C22	2.423	2.840	2.510	2.200	2.209	2.436	E4	10.80
9)	tC	C24	2.432	2.772	2.450	2.161	2.155	2.394	E4	10.62
10)	tC	C26	2.319	2.709	2.406	2.137	2.015	2.317	E4	11.53
11)	tC	C28	2.021	2.356	2.104	1.873	1.841	2.039	E4	10.16
12)	tC	C30	1.799	2.063	1.863	2.013	1.652	1.878	E4	8.83
13)	tC	C32	1.643	1.872	1.718	1.186	1.254	1.535	E4	19.53
14)	tC	C34	1.367	1.551	1.502	1.141	1.378	1.388	E4	11.46
15)	tC	C36	1.034	1.168	1.188	0.925	0.988	1.060	E4	10.76
16)	tC	C38	7.201	8.395	8.389	6.574	7.621	7.636	E3	10.27
17)	tC	C40	5.866	6.687	6.507	5.147	6.028	6.047	E3	10.00
18)	tC	c42	4.967	5.805	5.464	4.397	5.259	5.178	E3	10.29
19)	TC	Pristane	2.592	2.935	2.513	2.288	2.304	2.526	E4	10.43
20)	TC	Phytane	2.414	2.957	2.616	2.363	2.356	2.541	E4	10.06
21)	sC	o-terphenyl	2.709	3.183	2.823	2.559	2.556	2.766	E4	9.35
22)	tC	TPHC - total	2.282	2.624	2.440	2.600	2.736	2.536	E4	6.98

(#) = Out of Range

Evaluate Continuing Calibration Report

Data File : C:\HPCHEM\1\DATA\980327\T04706.D Vial: 2
Acq On : 27 Mar 98 3:16 pm Operator: DEINHARDT
Sample : 50 PPM STANDARD Inst : FID/TCD
Misc : Multiplr: 1.00
IntFile : TPHCINT.E

Method : C:\HPCHEM\1\METHODS\TPH27.M (Chemstation Integrator)
Title : TPHC Calibration 06/05/97 21 peaks
Last Update : Thu Mar 19 07:39:01 1998
Response via : Multiple Level Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
Max. RRF Dev : 25% Max. Rel. Area : 200%

		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
1	tC	C8	18.259	16.651 E3	8.8	94	0.01
2	tC	C10	19.534	18.921 E3	3.1	101	0.00
3	TC	C12	21.365	20.448 E3	4.3	101	0.00
4	tC	C14	22.088	20.910 E3	5.3	101	0.00
5	tC	C16	22.714	21.256 E3	6.4	101	0.00
6	tC	C18	25.919	23.759 E3	8.3	103	0.01
7	tC	C20	24.542	22.762 E3	7.3	102	0.01
8	tC	C22	24.364	22.827 E3	6.3	103	0.01
9	tC	C24	23.940	22.386 E3	6.5	102	0.01
10	tC	C26	23.170	22.081 E3	4.7	106	0.01
11	tC	C28	20.391	19.317 E3	5.3	109	0.01
12	tC	C30	18.781	17.288 E3	7.9	112	0.01
13	tC	C32	15.348	15.035 E3	2.0	107	0.01
14	tC	C34	13.879	14.158 E3	-2.0	120	0.02
15	tC	C36	10.605	11.317 E3	-6.7	125	0.02
16	tC	C38	7.636	8.520 E3	-11.6	126	0.03
17	tC	C40	6.047	6.617 E3	-9.4	125	0.03
18	tC	c42	5.178	5.405 E3	-4.4	122	0.05
19	TC	Pristane	25.262	22.814 E3	9.7	101	0.01
20	TC	Phytane	25.410	23.954 E3	5.7	102	0.01
21	sC	o-terphenyl	27.662	25.952 E3	6.2	103	0.01
22	tC	TPHC - total	25.364	22.924 E3	9.6	101	2.56#

(#) = Out of Range SPCC's out = 0 CCC's out = 0

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Report of Analysis
U.S. Army, Fort Monmouth Environmental Laboratory
NJDEP Certification # 13461

Surrogate Recovery Report

Lab. ID # : 3437

Location #: BLDG. 290

Sample		Surrogate Added (ppm)	Amount Recovered (ppm)	Percent Recovery
3437.02		10.00	10.38	**103.84**
3437.03		10.00	10.11	**101.09**
3437.04		10.00	10.70	**107.01**
3437.06		10.00	10.56	**105.56**
METHOD BLANK	27-Mar-98	10.00	10.31	**103.10**

Surrogate Added : **o-Terphenyl**

3/31/98

a

U.S. Army, Fort Monmouth Environmental Laboratory
NJDEP Certification # 13461

Matrix Spike Recovery Report

Lab. ID # : 3437

Location #: BLDG. 290

Sample	Spike Amount Added (ppm)	Sample Amount (ppm)	Matrix Spike Amount (ppm)	Percent Recovery	QC Limits %
3437.04MS	1000	53.85	1016.39	96.25	75-125
3437.04MSD	1000	53.85	1049.29	99.54	75-125

RPD	3.36	20.00

3/31/98

10

Report of Analysis
U.S. Army, Fort Monmouth Environmental Laboratory
NJDEP Certification # 13461

Blank Spike Recovery Report

Lab. ID # : 3437

Location #: BLDG. 290

Sample	Date Extracted	Spike Amount Added (ppm)	Matrix Spike Amount (ppm)	Percent Recovery	QC Limits %
Blank Spike	27-Mar-98	1000	1062.16	106.22	75-125

3/31/98

Quantitation Report (QT Reviewed)

Data File : C:\HPCHEM\1\DATA\980327\T04709.D Vial: 5
Acq On : 27 Mar 98 5:39 pm Operator: DEINHARDT
Sample : 3437.02 Inst : FID/TCD
Misc : Multiplr: 1.00
IntFile : TPHCINT.E
Quant Time: Mar 30 8:10 1998 Quant Results File: TPH27.RES

Quant Method : C:\HPCHEM\1\METHODS\TPH27.M (Chemstation Integrator)
Title : TPHC Calibration 06/05/97 21 peaks
Last Update : Thu Mar 19 07:39:01 1998
Response via : Initial Calibration
DataAcq Meth : TPH27.M

Volume Inj. : 1 ul
Signal Phase : HP-5
Signal Info : 30m x 0.32mm

Compound	R.T.	Response	Conc Units
System Monitoring Compounds			
21) sC o-terphenyl	13.95	287224	10.384 mg/L
Spiked Amount 10.000 Range 8 - 13		Recovery =	103.84%#
Target Compounds			
2) tC C10	8.82	1083	0.055 mg/L
3) TC C12	10.36	35514	1.662 mg/L
4) tC C14	11.52	1818	0.082 mg/L
5) tC C16	12.41	3407	0.150 mg/L
6) tC C18	12.99	1384	0.053 mg/L
7) tC C20	13.44	2963	0.121 mg/L
8) tC C22	14.15	3911	0.161 mg/L
9) tC C24	14.95	10982	0.459 mg/L
10) tC C26	15.72	2575	0.111 mg/L
11) tC C28	16.17	3597	0.176 mg/L
12) tC C30	17.01	1357	0.072 mg/L
13) tC C32	17.43	1705	0.111 mg/L
14) tC C34	17.83	1209	0.087 mg/L
15) tC C36	18.52	1229	0.116 mg/L
19) TC Pristane	12.99	1384	0.055 mg/L
20) TC Phytane	13.44	2963	0.117 mg/L
22) tC TPHC - total	13.95	2240429	88.331 mg/L m

(f)=RT Delta > 1/2 Window (m)=manual int.

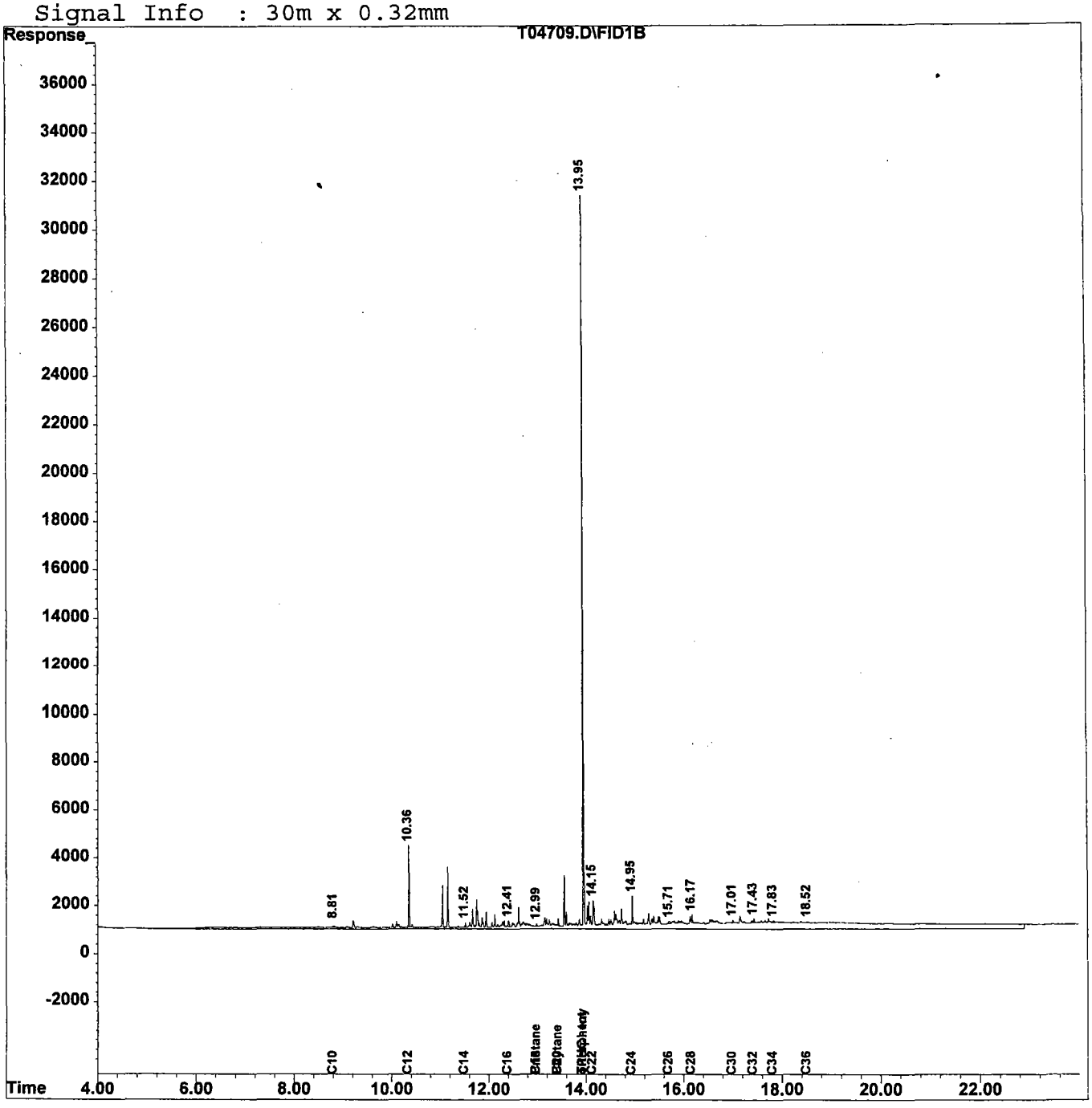
12

Quantitation Report

Data File : C:\HPCHEM\1\DATA\980327\T04709.D Vial: 5
Acq On : 27 Mar 98 5:39 pm Operator: DEINHARDT
Sample : 3437.02 Inst : FID/TCD
Misc : Multiplr: 1.00
IntFile : TPHCINT.E
Quant Time: Mar 30 8:10 1998 Quant Results File: TPH27.RES

Quant Method : C:\HPCHEM\1\METHODS\TPH27.M (Chemstation Integrator)
Title : TPHC Calibration 06/05/97 21 peaks
Last Update : Thu Mar 19 07:39:01 1998
Response via : Multiple Level Calibration
DataAcq Meth : TPH27.M

Volume Inj. : 1 ul
Signal Phase : HP-5
Signal Info : 30m x 0.32mm



Data File : C:\HPCHEM\1\DATA\980327\T04710.D Vial: 6
Acq On : 27 Mar 98 6:26 pm Operator: DEINHARDT
Sample : 3437.03 Inst : FID/TCD
Misc : Multiplr: 1.00
IntFile : TPHCINT.E
Quant Time: Mar 30 8:10 1998 Quant Results File: TPH27.RES

Quant Method : C:\HPCHEM\1\METHODS\TPH27.M (Chemstation Integrator)
Title : TPHC Calibration 06/05/97 21 peaks
Last Update : Thu Mar 19 07:39:01 1998
Response via : Initial Calibration
DataAcq Meth : TPH27.M

Volume Inj. : 1 ul
Signal Phase : HP-5
Signal Info : 30m x 0.32mm

Compound	R.T.	Response	Conc Units
System Monitoring Compounds			
21) sC o-terphenyl	13.95	279625	10.109 mg/L
Spiked Amount 10.000	Range 8 - 13	Recovery = 101.09%#	
Target Compounds			

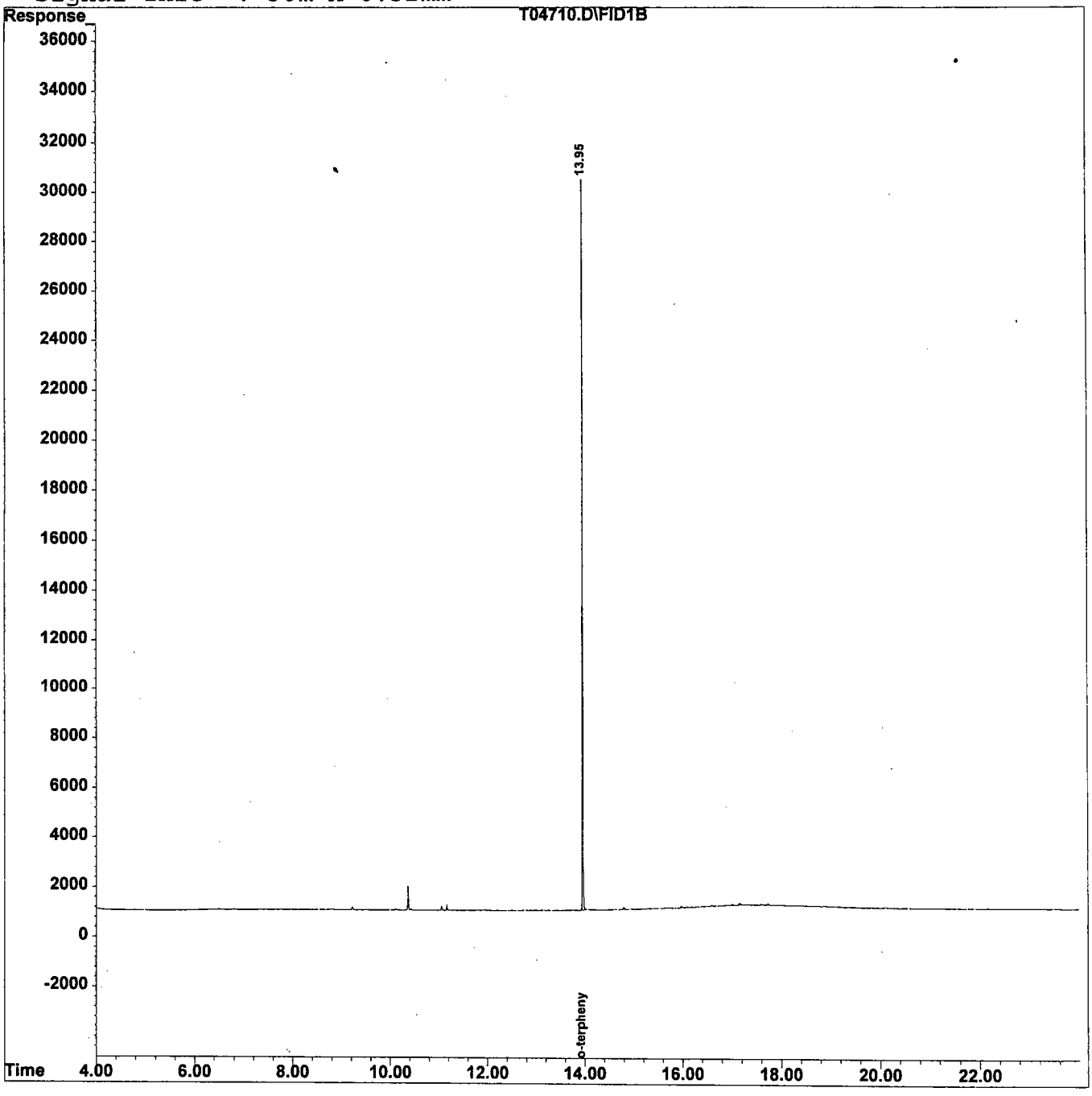
(f)=RT Delta > 1/2 Window (m)=manual int.

14

Data File : C:\HPCHEM\1\DATA\980327\T04710.D Vial: 6
Acq On : 27 Mar 98 6:26 pm Operator: DEINHARDT
Sample : 3437.03 Inst : FID/TCD
Misc : Multiplr: 1.00
IntFile : TPHCINT.E
Quant Time: Mar 30 8:10 1998 Quant Results File: TPH27.RES

Quant Method : C:\HPCHEM\1\METHODS\TPH27.M (Chemstation Integrator)
Title : TPHC Calibration 06/05/97 21 peaks
Last Update : Thu Mar 19 07:39:01 1998
Response via : Multiple Level Calibration
DataAcq Meth : TPH27.M

Volume Inj. : 1 ul
Signal Phase : HP-5
Signal Info : 30m x 0.32mm



Quantitation Report (QT Reviewed)

Data File : C:\HPCHEM\1\DATA\980327\T04711.D Vial: 7
Acq On : 27 Mar 98 7:11 pm Operator: DEINHARDT
Sample : 3437.04 Inst : FID/TCD
Misc : Multiplr: 1.00
IntFile : TPHCINT.E
Quant Time: Mar 30 8:11 1998 Quant Results File: TPH27.RES

Quant Method : C:\HPCHEM\1\METHODS\TPH27.M (Chemstation Integrator)
Title : TPHC Calibration 06/05/97 21 peaks
Last Update : Thu Mar 19 07:39:01 1998
Response via : Initial Calibration
DataAcq Meth : TPH27.M

Volume Inj. : 1 ul
Signal Phase : HP-5
Signal Info : 30m x 0.32mm

Compound	R.T.	Response	Conc	Units
System Monitoring Compounds				
21) sC o-terphenyl	13.95	296019	10.701	mg/L
Spiked Amount 10.000 Range 8 - 13		Recovery = 107.01%#		
Target Compounds				
11) tC C28	16.40	1460	0.072	mg/L
12) tC C30	16.88	1165	0.062	mg/L
13) tC C32	17.59	1126	0.073	mg/L
15) tC C36	18.74	2322	0.219	mg/L
22) tC TPHC - total	13.95	1365739	53.846	mg/L m

(f)=RT Delta > 1/2 Window (m)=manual int.

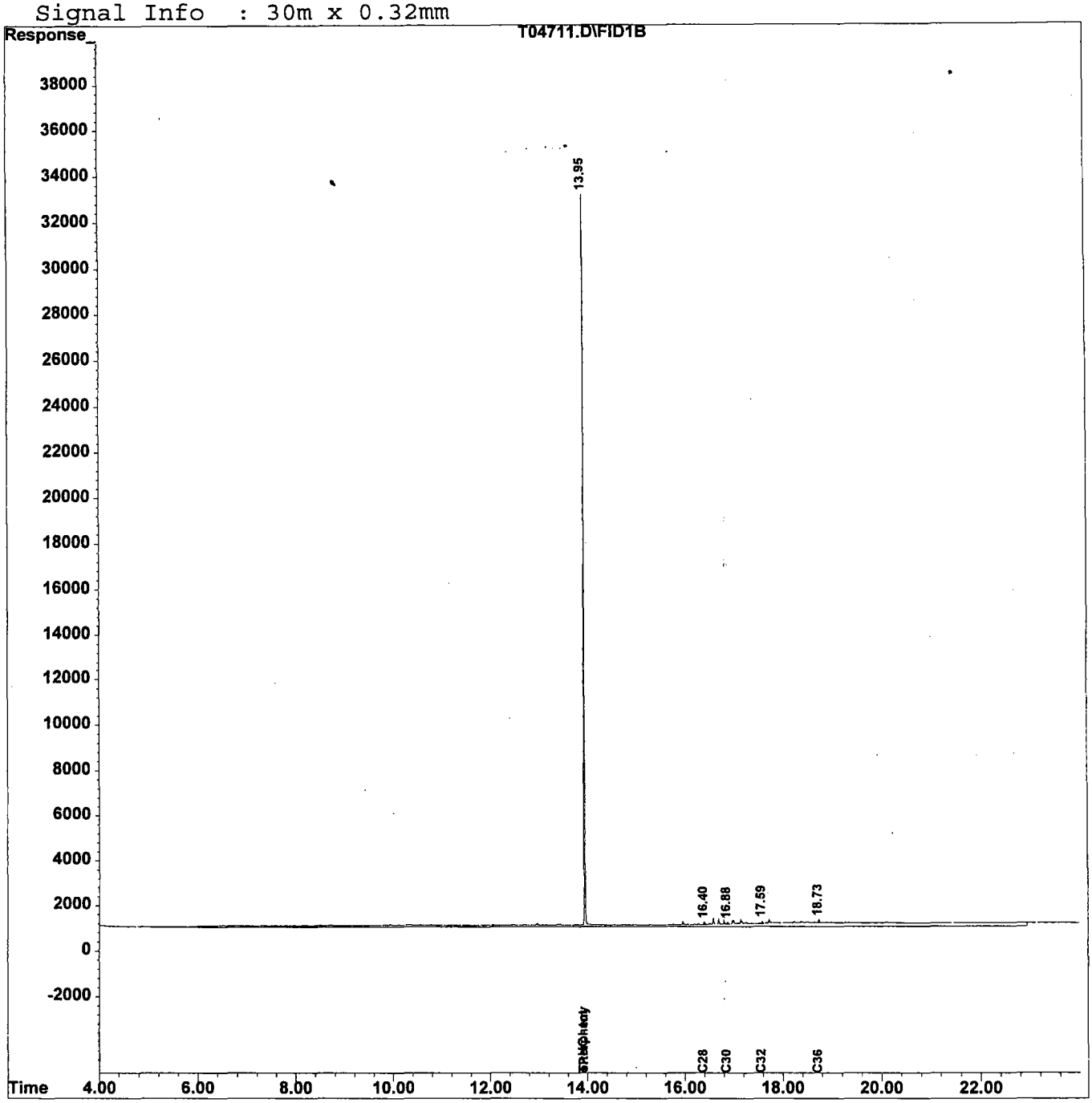
Quantitation Report

Data File : C:\HPCHEM\1\DATA\980327\T04711.D
Acq On : 27 Mar 98 7:11 pm
Sample : 3437.04
Misc :
IntFile : TPHCINT.E
Quant Time: Mar 30 8:11 1998 Quant Results File: TPH27.RES

Vial: 7
Operator: DEINHARDT
Inst : FID/TCD
Multiplr: 1.00

Quant Method : C:\HPCHEM\1\METHODS\TPH27.M (Chemstation Integrator)
Title : TPHC Calibration 06/05/97 21 peaks
Last Update : Thu Mar 19 07:39:01 1998
Response via : Multiple Level Calibration
DataAcq Meth : TPH27.M

Volume Inj. : 1 ul
Signal Phase : HP-5
Signal Info : 30m x 0.32mm



Quantitation Report (QT Reviewed)

Data File : C:\HPCHEM\1\DATA\980327\T04714.D Vial: 10
Acq On : 27 Mar 98 9:25 pm Operator: DEINHARDT
Sample : 3437.06 Inst : FID/TCD
Misc : Multiplr: 1.00
IntFile : TPHCINT.E
Quant Time: Mar 30 8:13 1998 Quant Results File: TPH27.RES

Quant Method : C:\HPCHEM\1\METHODS\TPH27.M (Chemstation Integrator)
Title : TPHC Calibration 06/05/97 21 peaks
Last Update : Thu Mar 19 07:39:01 1998
Response via : Initial Calibration
DataAcq Meth : TPH27.M

Volume Inj. : 1 ul
Signal Phase : HP-5
Signal Info : 30m x 0.32mm

Compound	R.T.	Response	Conc	Units
System Monitoring Compounds				
21) sC o-terphenyl	13.95	291994	10.556	mg/L
Spiked Amount 10.000	Range 8 - 13	Recovery = 105.56%#		
Target Compounds				

(f)=RT Delta > 1/2 Window (m)=manual int.

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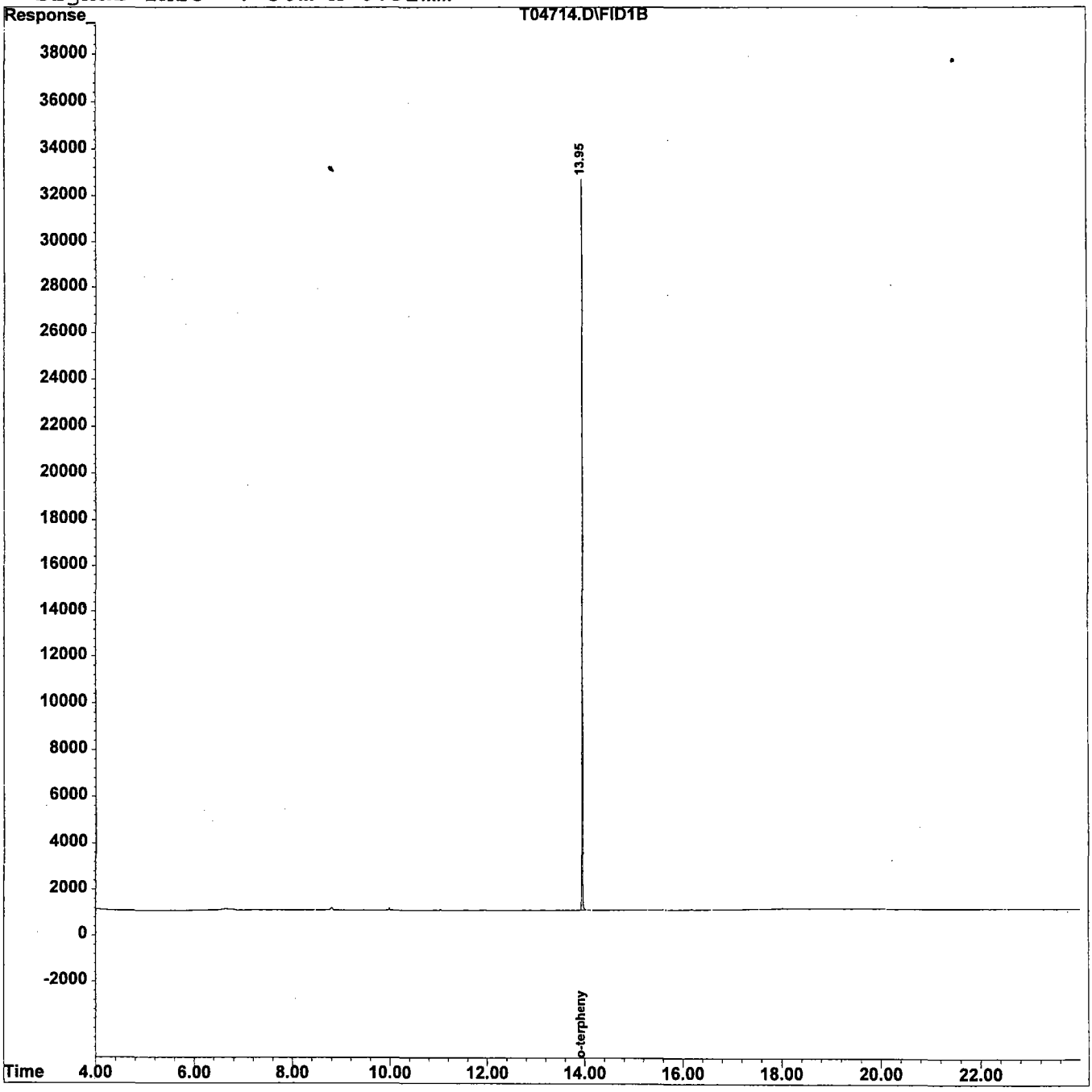
Quantitation Report

Data File : C:\HPCHEM\1\DATA\980327\T04714.D Vial: 10
Acq On : 27 Mar 98 9:25 pm Operator: DEINHARDT
Sample : 3437.06 Inst : FID/TCD
Misc : Multiplr: 1.00
IntFile : TPHCINT.E
Quant Time: Mar 30 8:13 1998 Quant Results File: TPH27.RES

Quant Method : C:\HPCHEM\1\METHODS\TPH27.M (Chemstation Integrator)
Title : TPHC Calibration 06/05/97 21 peaks
Last Update : Thu Mar 19 07:39:01 1998
Response via : Multiple Level Calibration
DataAcq Meth : TPH27.M

Volume Inj. : 1 ul
Signal Phase : HP-5
Signal Info : 30m x 0.32mm



19

LABORATORY DELIVERABLES CHECKLIST AND NON-CONFORMANCE SUMMARY

THIS FORM MUST BE COMPLETED BY THE LABORATORY OR ENVIRONMENTAL CONSULTANT AND ACCOMPANY ALL DATA SUBMISSIONS

The following Laboratory Deliverables checklist and Non-Conformance Summary shall be included in the data submission. All deviations from the accepted methodology and procedures, of performance values outside acceptable ranges shall be summarized in the Non-Conformance Summary. The Technical Requirements for Site Remediation, effective June 7, 1993, provides further details. The document shall be bound and paginated, contain a table of contents, and all pages shall be legible. Incomplete packages will be returned or held without review until the data package is completed.

It is recommended that the analytical results summary sheets listing all targeted and non-targeted compounds with the method detection limits, practical quantitation limits, and the laboratory and/or sample numbers be included in one section of the data package <u>and</u> in the main body of the report.

1. Cover page, Title Page listing Lab Certification #, facility name and address, & date of report submitted ✓
2. Table of Contents submitted ✓
3. Summary Sheets listing analytical results for all targeted and non-targeted compounds submitted ✓
4. Document paginated and legible ✓
5. Chain of Custody submitted ✓
6. Samples submitted to lab within 48 hours of sample collection ✓
7. Methodology Summary submitted ✓
8. Laboratory Chronicle and Holding Time Check submitted ✓
9. Results submitted on a dry weight basis ✓
10. Method Detection Limits submitted ✓
11. Lab certified by NJDEP for parameters of appropriate category of parameters or a member of the USEPA CLP ✓

Laboratory Manager or Environmental Consultant's Signature _____

Date 4/6/98

Laboratory Certification #13461

*Refer to NJAC 7:26E - Appendix A, Section IV - Reduced Data Deliverables - Non-USEPA/CLP Methods for further guidance

US ARMY FT. MONMOUTH ENVIRONMENTAL LABORATORY
NJDEPE # 13461

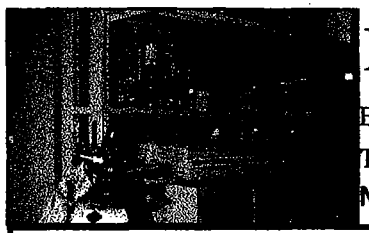
REPORT OF ANALYSIS

Client: U.S. Army
DPW, SELFM-PW-EV
Bldg. 173
Ft. Monmouth, NJ 07703

Project: Volatiles - EPA Method 624
Bldg. 290

Project # 3437
Date Rec. 03/26/98
Date Compl. 03/30/98
Released by:

Daniel K. Wright
Laboratory Director



Fort Monmouth Environmental Testing Laboratory

Bldg. 173, SELFM-PW-EV, Fort Monmouth, NJ 07703
Tel (732)532-4359 Fax (732)532-3484 EMail:appleby@doim6.monmouth.army.mil
NJDEP Certification #13461

Chain of Custody Record

Customer:

Phone #:

( )DERA ( )OMA ( )Other:_____

Project No:

Location: Bldg-290

Samplers Name / Company :

Comments: OVA #2-AS1903

Lab Sample I.D.	Sample Location	Date	Time	Sample Type	# bottles	VOA + 15	TPH						Remarks / Preservation Method
3437.01	TRIP BLANK	3-26-98	~~1348~~	AQ	2	X							HCL
02	290-B-1		1348	SOIL	1		X						9 PPM) 24-30" (ALC)
03	290-B-2		1433	SOIL	1		X						100 PPM) 22-26" (ML)
04	290-B-3		1507	SOIL	1		X						90 PPM)
(MW) ~~05~~	~~290-B-4~~			~~SOIL~~	~~1~~		~~X~~						
05	290-B-2-W		1550	AQ	2	X							
06	290-B-4		1625	SOIL	1		X						

Relinquished by (signature):	Date/Time:	Received by (signature):	Relinquished by (signature):	Date/Time:	Received by (signature):
	3-26-98 1640				
Relinquished by (signature):	Date/Time:	Received by (signature):	Relinquished by (signature):	Date/Time:	Received by (signature):

Report Type: (_)Full, (_)Reduced, (_)Standard, (_)Screen / non-certified

Turnaround time: (_)Standard 4 wks, (_)Rush ____ Days, (_)ASAP Verbal ____ Hrs.

Remarks:

print legibly

~me1a23.xls2/20/98

Volatile Analysis Report
U.S. Army, Fort Monmouth Environmental Laboratory
NJDEP Certification #13461

Data File Name	**V03392.D**	Sample Name	**VBLK41**
Operator	**Skelton**	Field ID	**VBLK41**
Date Acquired	**30 Mar 1998 20:22**	Sample Multiplier	**1**

CAS#	Compound Name	R.T.	Response	Result	GW Criteria	MDL	Qualifiers
107028	Acrolein			not detected	nle	6.25 ug/L	
107131	Acrylonitrile			not detected	nle	6.25 ug/L	
75650	tert-Butyl alcohol			not detected	nle	12.50 ug/L	
1634044	Methyl-tert-Butyl ether			not detected	nle	2.50 ug/L	
108203	Di-isopropyl ether			not detected	nle	1.25 ug/L	
	Dichlorodifluoromethane			not detected	nle	3.63 ug/L	
74-87-3	Chloromethane			not detected	30	0.79 ug/L	
75-01-4	Vinyl Chloride			not detected	5	2.61 ug/L	
74-83-9	Bromomethane			not detected	10	1.45 ug/L	
75-00-3	Chloroethane			not detected	nle	2.20 ug/L	
75-69-4	Trichlorofluoromethane			not detected	nle	1.31 ug/L	
75-35-4	1,1-Dichloroethene			not detected	2	0.74 ug/L	
67-64-1	Acetone			not detected	700	1.57 ug/L	
75-15-0	Carbon Disulfide			not detected	nle	0.54 ug/L	
75-09-2	Methylene Chloride			not detected	2	1.66 ug/L	
156-60-5	trans-1,2-Dichloroethene			not detected	100	0.50 ug/L	
75-35-3	1,1-Dichloroethane			not detected	70	0.83 ug/L	
108-05-4	Vinyl Acetate			not detected	nle	2.07 ug/L	
78-93-3	2-Butanone			not detected	300	2.06 ug/L	
	cis-1,2-Dichloroethene			not detected	10	0.65 ug/L	
67-66-3	Chloroform			not detected	6	0.43 ug/L	
75-55-6	1,1,1-Trichloroethane			not detected	30	0.81 ug/L	
56-23-5	Carbon Tetrachloride			not detected	2	1.20 ug/L	
71-43-2	Benzene			not detected	1	0.51 ug/L	
107-06-2	1,2-Dichloroethane			not detected	2	1.27 ug/L	
79-01-6	Trichloroethene			not detected	1	0.94 ug/L	
78-87-5	1,2-Dichloropropane			not detected	1	0.78 ug/L	
75-27-4	Bromodichloromethane			not detected	1	0.77 ug/L	
110-75-8	2-Chloroethyl vinyl ether			not detected	nle	1.05 ug/L	
10061-01-5	cis-1,3-Dichloropropene			not detected	nle	0.60 ug/L	
108-10-1	4-Methyl-2-Pentanone			not detected	400	1.33 ug/L	
108-88-3	Toluene			not detected	1000	0.73 ug/L	
10061-02-6	trans-1,3-Dichloropropene			not detected	nle	1.43 ug/L	
79-00-5	1,1,2-Trichloroethane			not detected	3	1.49 ug/L	
127-18-4	Tetrachloroethene			not detected	1	0.92 ug/L	
591-78-6	2-Hexanone			not detected	nle	1.12 ug/L	
126-48-1	Dibromochloromethane			not detected	10	1.36 ug/L	
108-90-7	Chlorobenzene			not detected	4	0.66 ug/L	
100-41-4	Ethylbenzene			not detected	700	1.14 ug/L	
1330-20-7	m+p-Xylenes			not detected	nle	2.53 ug/L	
1330-20-7	o-Xylene			not detected	nle	1.92 ug/L	
100-42-5	Styrene			not detected	100	1.57 ug/L	
75-25-2	Bromoform			not detected	4	1.68 ug/L	
79-34-5	1,1,2,2-Tetrachloroethane			not detected	2	1.71 ug/L	
541-73-1	1,3-Dichlorobenzene			not detected	600	2.51 ug/L	
106-46-7	1,4-Dichlorobenzene			not detected	74	3.08 ug/L	
95-50-1	1,2-Dichlorobenzene			not detected	600	2.75 ug/L	

Qualifiers

B = Compound found in related blank
E = Value above linear range
D = Value from dilution
MDL = Method Detection Limit
NLE = No Limit Established
R.T. = Retention Time

1E

VOLATILE ORGANICS ANALYSIS DATA SHEET

TENTATIVELY IDENTIFIED COMPOUNDS

FIELD ID.

VBLK41

Lab Name: FMETL NJDEP # 13461

Project: Case No.: 3437 Location: B.290 SDG No.:

Matrix: (soil/water) WATER Lab Sample ID: VBLK41

Sample wt/vol: 5.0 (g/ml) ML Lab File ID: V03392.D

Level: (low/med) LOW Date Received: 03/26/98

% Moisture: not dec. Date Analyzed: 03/30/98

GC Column: Rtx502.2 ID: 0.25 (mm) Dilution Factor: 1.0

Soil Extract Volume: (uL) Soil Aliquot Volume: (uL)

CONCENTRATION UNITS:

(ug/L or ug/Kg) UG/L

Number TICs found: 1

CAS NO.	COMPOUND	RT	EST. CONC.	Q
1.	unknown	12.16	27	J

Volatile Analysis Report
U.S. Army, Fort Monmouth Environmental Laboratory
NJDEP Certification #13461

Data File Name	**v03408.d**	Sample Name	**3437.01**
Operator	**Skelton**	Field ID	**Trip Blank**
Date Acquired	**31 Mar 1998 8:57**	Sample Multiplier	**1**

CAS#	Compound Name	R.T.	Response	Result	GW Criteria	MDL	Qualifiers
107028	Acrolein			not detected	nle	6.25 ug/L	
107131	Acrylonitrile			not detected	nle	6.25 ug/L	
75650	tert-Butyl alcohol			not detected	nle	12.50 ug/L	
1634044	Methyl-tert-Butyl ether			not detected	nle	2.50 ug/L	
108203	Di-isopropyl ether			not detected	nle	1.25 ug/L	
	Dichlorodifluoromethane			not detected	nle	3.63 ug/L	
74-87-3	Chloromethane			not detected	30	0.79 ug/L	
75-01-4	Vinyl Chloride			not detected	5	2.61 ug/L	
74-83-9	Bromomethane			not detected	10	1.45 ug/L	
75-00-3	Chloroethane			not detected	nle	2.20 ug/L	
75-69-4	Trichlorofluoromethane			not detected	nle	1.31 ug/L	
75-35-4	1,1-Dichloroethene			not detected	2	0.74 ug/L	
67-64-1	Acetone			not detected	700	1.57 ug/L	
75-15-0	Carbon Disulfide			not detected	nle	0.54 ug/L	
75-09-2	Methylene Chloride			not detected	2	1.66 ug/L	
156-60-5	trans-1,2-Dichloroethene			not detected	100	0.50 ug/L	
75-35-3	1,1-Dichloroethane			not detected	70	0.83 ug/L	
108-05-4	Vinyl Acetate			not detected	nle	2.07 ug/L	
78-93-3	2-Butanone			not detected	300	2.06 ug/L	
	cis-1,2-Dichloroethene			not detected	10	0.65 ug/L	
67-66-3	Chloroform			not detected	6	0.43 ug/L	
75-55-6	1,1,1-Trichloroethane			not detected	30	0.81 ug/L	
56-23-5	Carbon Tetrachloride			not detected	2	1.20 ug/L	
71-43-2	Benzene			not detected	1	0.51 ug/L	
107-06-2	1,2-Dichloroethane			not detected	2	1.27 ug/L	
79-01-6	Trichloroethene			not detected	1	0.94 ug/L	
78-87-5	1,2-Dichloropropane			not detected	1	0.78 ug/L	
75-27-4	Bromodichloromethane			not detected	1	0.77 ug/L	
110-75-8	2-Chloroethyl vinyl ether			not detected	nle	1.05 ug/L	
10061-01-5	cis-1,3-Dichloropropene			not detected	nle	0.60 ug/L	
108-10-1	4-Methyl-2-Pentanone			not detected	400	1.33 ug/L	
108-88-3	Toluene			not detected	1000	0.73 ug/L	
10061-02-6	trans-1,3-Dichloropropene			not detected	nle	1.43 ug/L	
79-00-5	1,1,2-Trichloroethane			not detected	3	1.49 ug/L	
127-18-4	Tetrachloroethene			not detected	1	0.92 ug/L	
591-78-6	2-Hexanone			not detected	nle	1.12 ug/L	
126-48-1	Dibromochloromethane			not detected	10	1.36 ug/L	
108-90-7	Chlorobenzene			not detected	4	0.66 ug/L	
100-41-4	Ethylbenzene			not detected	700	1.14 ug/L	
1330-20-7	m+p-Xylenes			not detected	nle	2.53 ug/L	
1330-20-7	o-Xylene			not detected	nle	1.92 ug/L	
100-42-5	Styrene			not detected	100	1.57 ug/L	
75-25-2	Bromoform			not detected	4	1.68 ug/L	
79-34-5	1,1,2,2-Tetrachloroethane			not detected	2	1.71 ug/L	
541-73-1	1,3-Dichlorobenzene			not detected	600	2.51 ug/L	
106-46-7	1,4-Dichlorobenzene			not detected	74	3.08 ug/L	
95-50-1	1,2-Dichlorobenzene			not detected	600	2.75 ug/L	

Qualifiers

B = Compound found in related blank
E = Value above linear range
D = Value from dilution
MDL = Method Detection Limit
NLE = No Limit Established
R.T. = Retention Time

1E

VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

FIELD ID.

Trip Blank

Lab Name: FMETL NJDEP # 13461

Project: Case No.: 3437 Location: B.290 SDG No.:

Matrix: (soil/water) WATER Lab Sample ID: 3437.01

Sample wt/vol: 5.0 (g/ml) ML Lab File ID: V03408.D

Level: (low/med) LOW Date Received: 03/26/98

% Moisture: not dec. Date Analyzed: 03/31/98

GC Column: Rtx502.2 ID: 0.25 (mm) Dilution Factor: 1.0

Soil Extract Volume: (uL) Soil Aliquot Volume: (uL)

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/L

Number TICs found: 0

CAS NO.	COMPOUND	RT	EST. CONC.	Q

Volatile Analysis Report
U.S. Army, Fort Monmouth Environmental Laboratory
NJDEP Certification #13461

Data File Name	**v03409.d**	Sample Name	**3437.05**
Operator	**Skelton**	Field ID	**290-B-2-W**
Date Acquired	**31 Mar 1998 9:41**	Sample Multiplier	**1**

CAS#	Compound Name	R.T.	Response	Result	GW Criteria	MDL	Qualifiers
107028	Acrolein			not detected	nle	6.25 ug/L	
107131	Acrylonitrile			not detected	nle	6.25 ug/L	
75650	tert-Butyl alcohol			not detected	nle	12.50 ug/L	
1634044	Methyl-tert-Butyl ether			not detected	nle	2.50 ug/L	
108203	Di-isopropyl ether			not detected	nle	1.25 ug/L	
	Dichlorodifluoromethane			not detected	nle	3.63 ug/L	
74-87-3	Chloromethane			not detected	30	0.79 ug/L	
75-01-4	Vinyl Chloride			not detected	5	2.61 ug/L	
74-83-9	Bromomethane			not detected	10	1.45 ug/L	
75-00-3	Chloroethane			not detected	nle	2.20 ug/L	
75-69-4	Trichlorofluoromethane			not detected	nle	1.31 ug/L	
75-35-4	1,1-Dichloroethene			not detected	2	0.74 ug/L	
67-64-1	Acetone			not detected	700	1.57 ug/L	
75-15-0	Carbon Disulfide			not detected	nle	0.54 ug/L	
75-09-2	Methylene Chloride			not detected	2	1.66 ug/L	
156-60-5	trans-1,2-Dichloroethene			not detected	100	0.50 ug/L	
75-35-3	1,1-Dichloroethane			not detected	70	0.83 ug/L	
108-05-4	Vinyl Acetate			not detected	nle	2.07 ug/L	
78-93-3	2-Butanone			not detected	300	2.06 ug/L	
	cis-1,2-Dichloroethene			not detected	10	0.65 ug/L	
67-66-3	Chloroform			not detected	6	0.43 ug/L	
75-55-6	1,1,1-Trichloroethane			not detected	30	0.81 ug/L	
56-23-5	Carbon Tetrachloride			not detected	2	1.20 ug/L	
71-43-2	Benzene			not detected	1	0.51 ug/L	
107-06-2	1,2-Dichloroethane			not detected	2	1.27 ug/L	
79-01-6	Trichloroethene			not detected	1	0.94 ug/L	
78-87-5	1,2-Dichloropropane			not detected	1	0.78 ug/L	
75-27-4	Bromodichloromethane			not detected	1	0.77 ug/L	
110-75-8	2-Chloroethyl vinyl ether			not detected	nle	1.05 ug/L	
10061-01-5	cis-1,3-Dichloropropene			not detected	nle	0.60 ug/L	
108-10-1	4-Methyl-2-Pentanone			not detected	400	1.33 ug/L	
108-88-3	Toluene			not detected	1000	0.73 ug/L	
10061-02-6	trans-1,3-Dichloropropene			not detected	nle	1.43 ug/L	
79-00-5	1,1,2-Trichloroethane			not detected	3	1.49 ug/L	
127-18-4	Tetrachloroethene			not detected	1	0.92 ug/L	
591-78-6	2-Hexanone			not detected	nle	1.12 ug/L	
126-48-1	Dibromochloromethane			not detected	10	1.36 ug/L	
108-90-7	Chlorobenzene			not detected	4	0.66 ug/L	
100-41-4	Ethylbenzene			not detected	700	1.14 ug/L	
1330-20-7	m+p-Xylenes			not detected	nle	2.53 ug/L	
1330-20-7	o-Xylene			not detected	nle	1.92 ug/L	
100-42-5	Styrene			not detected	100	1.57 ug/L	
75-25-2	Bromoform			not detected	4	1.68 ug/L	
79-34-5	1,1,2,2-Tetrachloroethane			not detected	2	1.71 ug/L	
541-73-1	1,3-Dichlorobenzene			not detected	600	2.51 ug/L	
106-46-7	1,4-Dichlorobenzene			not detected	74	3.08 ug/L	
95-50-1	1,2-Dichlorobenzene			not detected	600	2.75 ug/L	

Qualifiers

B = Compound found in related blank
E = Value above linear range
D = Value from dilution
MDL = Method Detection Limit
NLE = No Limit Established
R.T. = Retention Time

1E

VOLATILE ORGANICS ANALYSIS DATA SHEET FIELD ID.

TENTATIVELY IDENTIFIED COMPOUNDS

B-4

Lab Name: FMETL NJDEP # 13461

Project: Case No.: 3437 Location: B.290 SDG No.:

Matrix: (soil/water) WATER Lab Sample ID: 3437.05

Sample wt/vol: 5.0 (g/ml) ML Lab File ID: V03409.D

Level: (low/med) LOW Date Received: 03/26/98

% Moisture: not dec. Date Analyzed: 03/31/98

GC Column: Rtx502.2 ID: 0.25 (mm) Dilution Factor: 1.0

Soil Extract Volume: (uL) Soil Aliquot Volume: (uL)

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/L

Number TICs found: 2

CAS NO.	COMPOUND	RT	EST. CONC.	Q
1. 001070-71-9	Propiolonitrile	5.16	5	JN
2. 000593-75-9	Methane, isocyano-	12.17	9	JN

Quantitation Report (QT Reviewed)

Data File : C:\HPCHEM\1\DATA\980330\V03392.D Vial: 7
Acq On : 30 Mar 1998 20:22 Operator: Skelton
Sample : VBLK41 Inst : GC/MS Ins
Misc : VBLK41 Multiplr: 1.00
MS Integration Params: gases.p
Quant Time: Mar 31 9:10 1998 Quant Results File: M62431.RES

Quant Method : C:\HPCHEM\1\METHODS\M62431.M (RTE Integrator)
Title : Volatile Organics by GC/MS Method 624/8240/TCLP
Last Update : Tue Mar 31 09:08:52 1998
Response via : Initial Calibration
DataAcq Meth : M62429

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Bromochloromethane	18.77	128	31942	30.00	ug/L	0.00
26) 1,4-Difluorobenzene	21.36	114	220384	30.00	ug/L	0.00
37) Chlorobenzene-d5	29.19	119	64135	30.00	ug/L	0.00
System Monitoring Compounds						
25) 1,2-Dichloroethane-d4	20.32	65	99897	30.31	ug/L	0.00
Spiked Amount 30.000	Range 76 - 114		Recovery = 101.03%			
35) Toluene-d8	25.35	98	272267	30.08	ug/L	0.00
Spiked Amount 30.000	Range 88 - 110		Recovery = 100.27%			
49) Bromofluorobenzene	32.19	95	119248	29.46	ug/L	0.00
Spiked Amount 30.000	Range 86 - 115		Recovery = 98.20%			

Target Compounds Qvalue

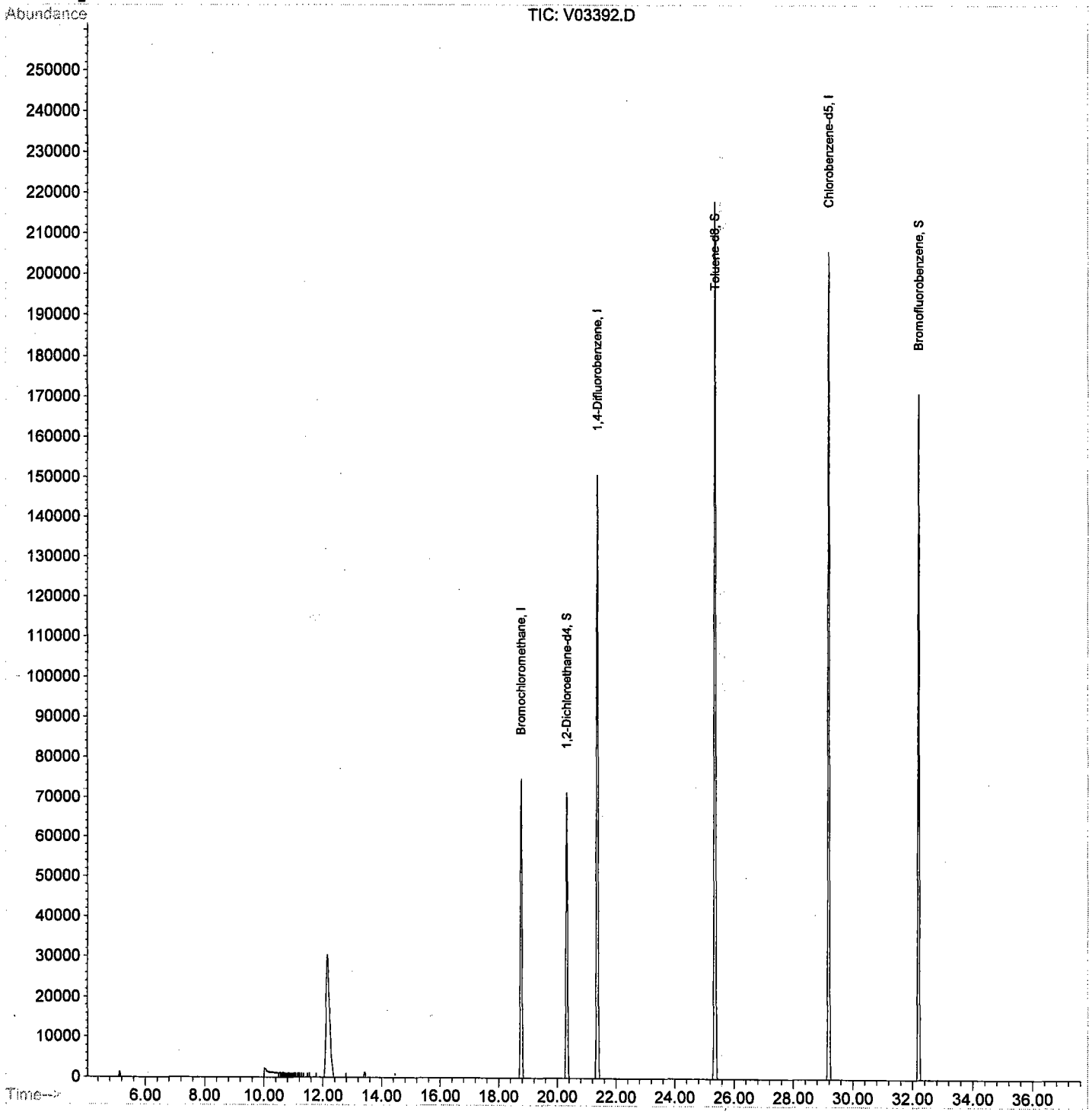
(#) = qualifier out of range (m) = manual integration

Quantitation Report

Data File : C:\HPCHEM\1\DATA\980330\V03392.D
Acq On : 30 Mar 1998 20:22
Sample : VBLK41
Misc : VBLK41
MS Integration Params: gases.p
Quant Time: Mar 31 9:10 1998

Vial: 7
Operator: Skelton
Inst : GC/MS Ins
Multiplr: 1.00
Quant Results File: M62431.RES

Method : C:\HPCHEM\1\METHODS\M62433.M (RTE Integrator)
Title : Volatile Organics by GC/MS Method 624/8240/TCLP
Last Update : Wed Apr 22 07:15:47 1998
Response via : Initial Calibration



Quantitation Report (QT Reviewed)

Data File : C:\HPCHEM\1\DATA\980330\V03408.D Vial: 8
Acq On : 31 Mar 1998 8:57 Operator: Skelton
Sample : 3437.01 Inst : GC/MS Ins
Misc : Trip Blank Multiplr: 1.00
MS Integration Params: gases.p
Quant Time: Apr 24 8:19 1998 Quant Results File: M62431.RES

Quant Method : C:\HPCHEM\1\METHODS\M62431.M (RTE Integrator)
Title : Volatile Organics by GC/MS Method 624/8240/TCLP
Last Update : Tue Mar 31 09:08:52 1998
Response via : Initial Calibration
DataAcq Meth : M62431

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Bromochloromethane	18.78	128	29377	30.00	ug/L	0.00
26) 1,4-Difluorobenzene	21.37	114	218649	30.00	ug/L	0.00
37) Chlorobenzene-d5	29.20	119	64278	30.00	ug/L	0.01
System Monitoring Compounds						
25) 1,2-Dichloroethane-d4	20.33	65	92533	30.53	ug/L	0.00
Spiked Amount 30.000	Range 76 - 114		Recovery = 101.77%			
35) Toluene-d8	25.37	98	272715	30.36	ug/L	0.01
Spiked Amount 30.000	Range 88 - 110		Recovery = 101.20%			
49) Bromofluorobenzene	32.21	95	117188	28.89	ug/L	0.01
Spiked Amount 30.000	Range 86 - 115		Recovery = 96.30%			

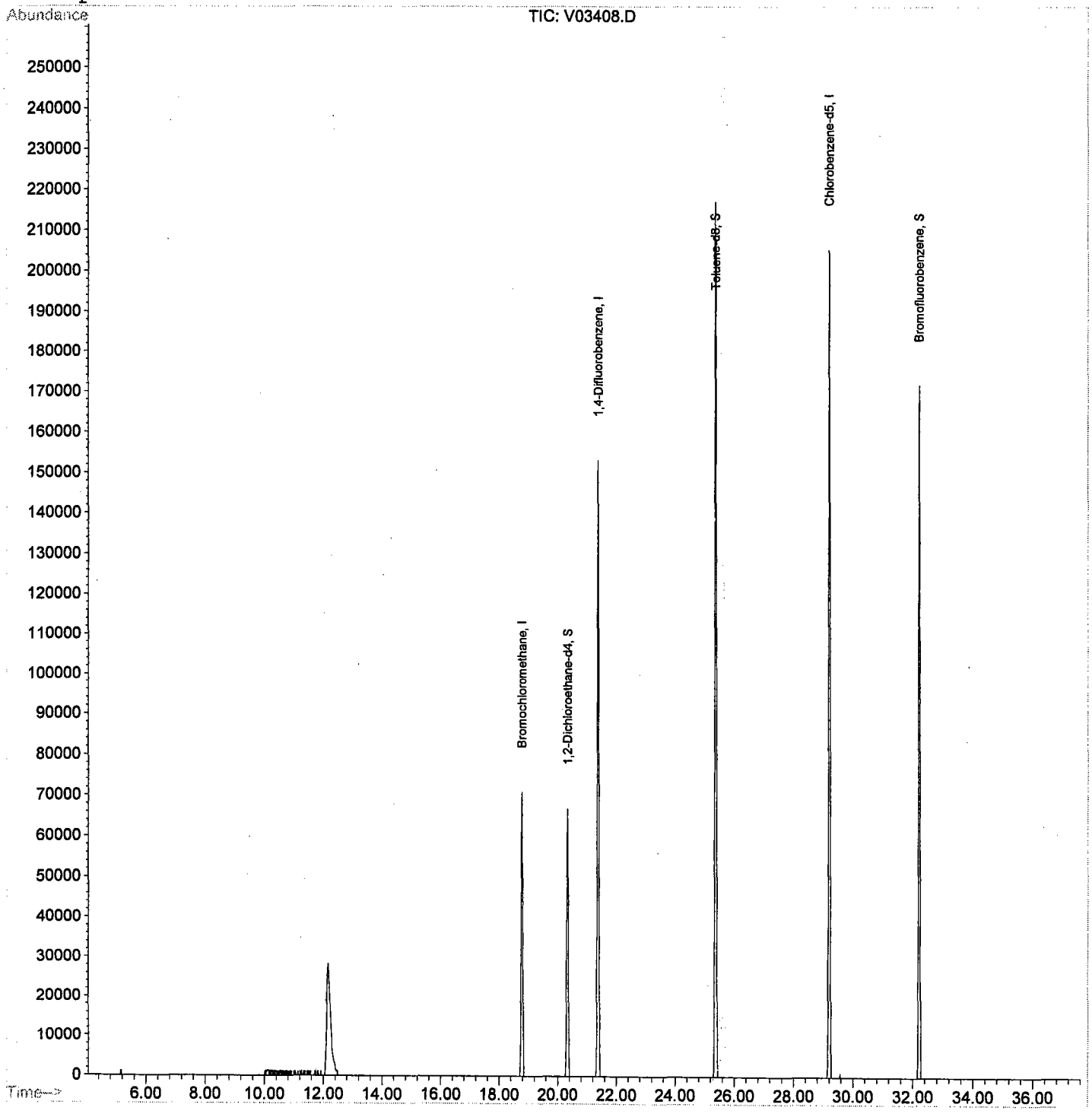
Target Compounds Qvalue

(#) = qualifier out of range (m) = manual integration

Quantitation Report

Data File : C:\HPCHEM\1\DATA\980330\V03408.D Vial: 8
Acq On : 31 Mar 1998 8:57 Operator: Skelton
Sample : 3437.01 Inst : GC/MS Ins
Misc : Trip Blank Multiplr: 1.00
MS Integration Params: gases.p
Quant Time: Apr 24 8:19 1998 Quant Results File: M62431.RES

Method : C:\HPCHEM\1\METHODS\M62433.M (RTE Integrator)
Title : Volatile Organics by GC/MS Method 624/8240/TCLP
Last Update : Wed Apr 22 07:15:47 1998
Response via : Initial Calibration



Quantitation Report (QT Reviewed)

Data File : C:\HPCHEM\1\DATA\980330\V03409.D Vial: 9
Acq On : 31 Mar 1998 9:41 Operator: Skelton
Sample : 3437.05 Inst : GC/MS Ins
Misc : 290-B-2-W Multiplr: 1.00
MS Integration Params: gases.p
Quant Time: Apr 24 8:21 1998 Quant Results File: M62431.RES

Quant Method : C:\HPCHEM\1\METHODS\M62431.M (RTE Integrator)
Title : Volatile Organics by GC/MS Method 624/8240/TCLP
Last Update : Tue Mar 31 09:08:52 1998
Response via : Initial Calibration
DataAcq Meth : M62431

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Bromochloromethane	18.78	128	27977	30.00	ug/L	0.01
26) 1,4-Difluorobenzene	21.38	114	202794	30.00	ug/L	0.01
37) Chlorobenzene-d5	29.21	119	60653	30.00	ug/L	0.02
System Monitoring Compounds						
25) 1,2-Dichloroethane-d4	20.33	65	87776	30.41	ug/L	0.01
Spiked Amount 30.000	Range 76 - 114		Recovery = 101.37%			
35) Toluene-d8	25.37	98	252003	30.25	ug/L	0.01
Spiked Amount 30.000	Range 88 - 110		Recovery = 100.83%			
49) Bromofluorobenzene	32.21	95	113388	29.62	ug/L	0.01
Spiked Amount 30.000	Range 86 - 115		Recovery = 98.73%			

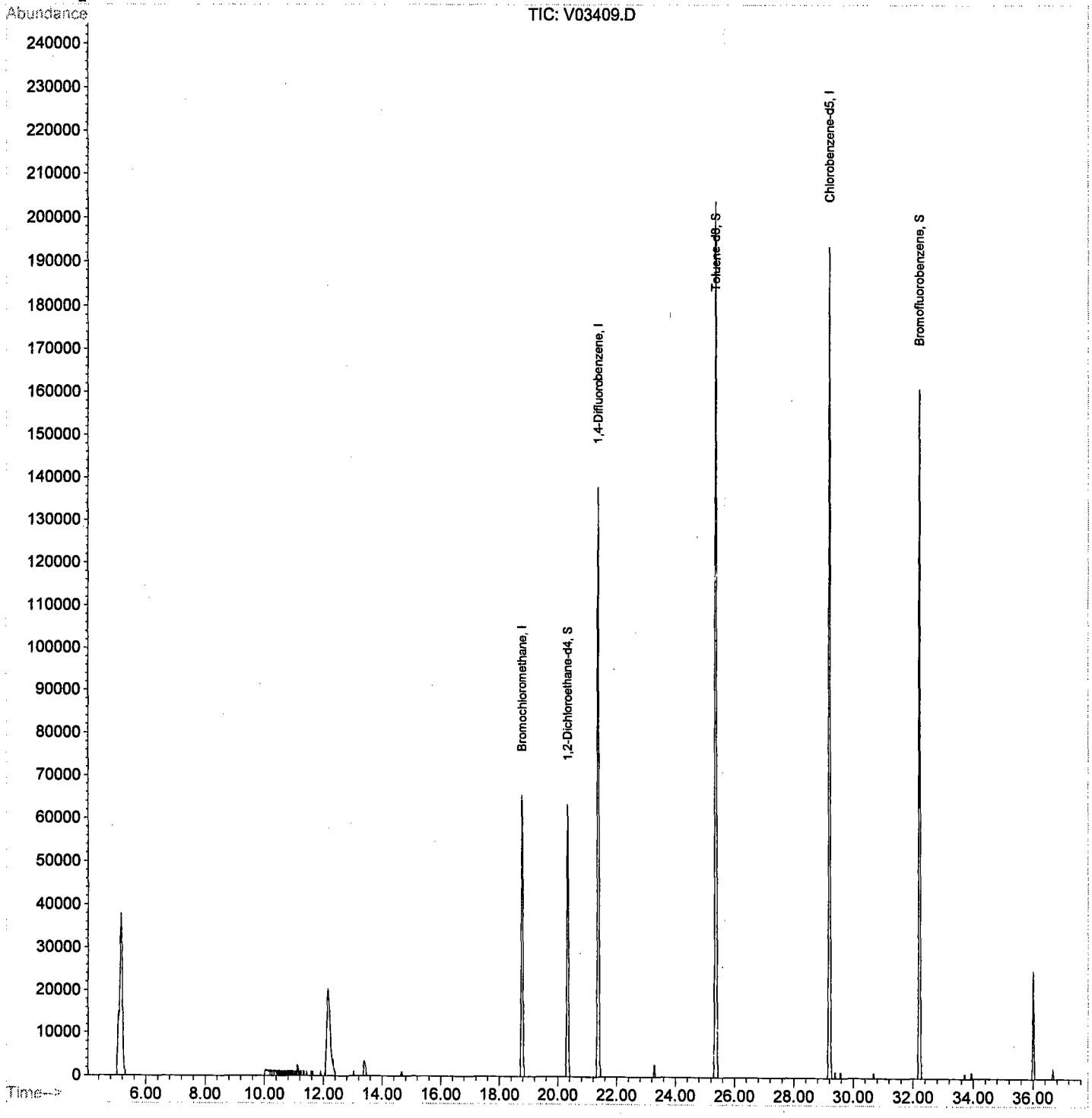
Target Compounds Qvalue

(#) = qualifier out of range (m) = manual integration

Quantitation Report

Data File : C:\HPCHEM\1\DATA\980330\V03409.D Vial: 9
Acq On : 31 Mar 1998 9:41 Operator: Skelton
Sample : 3437.05 Inst : GC/MS Ins
Misc : 290-B-2-W Multiplr: 1.00
MS Integration Params: gases.p
Quant Time: Apr 24 8:21 1998 Quant Results File: M62431.RES

Method : C:\HPCHEM\1\METHODS\M62433.M (RTE Integrator)
Title : Volatile Organics by GC/MS Method 624/8240/TCLP
Last Update : Wed Apr 22 07:15:47 1998
Response via : Initial Calibration



APPENDIX G

GROUNDWATER ANALYTICAL DATA PACKAGE

Project No.: 9421415
Log in No. : 22509E
P.O. No. : Pending
Date : 2/21/95
SDG No. : Army 1
NJDEPE Case #:93-11-30-1246-27

ANALYTICAL DATA REPORT
PACKAGE FOR

Aguilar Associates

30 Freneau Avenue

Matawan, NJ 07747

ATTN: Darryl Schmitt
REF: US Army Fort Monmouth, Well# and NJPDEPE Reg# 1-2930961
Sample Location Bldg. 290

LABORATORY NUMBER	SAMPLE IDENTIFICATION	TYPE OF SAMPLE

S E E N E X T P A G E

WE CERTIFY THAT THIS REPORT IS A TRUE REPORT OF RESULTS OBTAINED FROM OUR TESTS OF THIS MATERIAL.

RESPECTFULLY SUBMITTED,
NYTEST ENVIRONMENTAL INC.

NYS Lab ID. #10195
NJ Cert. #73469

REMO GIGANTE
EXEC. VICE PRESIDENT

NYTEST ENVIRONMENTAL Inc

LABORATORY NUMBER	SAMPLE IDENTIFICATION	WELL #	TYPE OF SAMPLE
2250912	290-1	1-2930961	Water
2250904	108-FB	-	Water
2250916	2562-TB	-	Water

Table of Contents

BLDG.#: 290 MW#: 1 NJDEPE WELL ID # 2930961

U.S. ARMY FORT MONMOUTH

MONITORING WELL SAMPLING DATASHEET

DATE: 11/8/91

IJO#94-0843 A

SAMPLING CONTRACTOR: Aguilar Associates Inc.

LABORATORY: NYTEST Environmental Inc. CERT #:73469

SAMPLERS NAMES: D. Schmith, C. Aguilar, N. Vonasek, E. Romero, S. Panizzi

WEATHER CONDITIONS: 58°F Sunny, windy

ELEVATION OF CASING SURVEY MARK: 13.90

TOTAL DEPTH OF WELL FROM TOP OF SURVEYORS MARK: 16.52 FT

DEPTH FROM SURVEYORS MARK TO SCREEN: 2.0 FT

LENGTH OF SCREENED SECTION: 10.5 FT.

DEPTH TO WATER PRIOR TO PURGING AND SAMPLING: 8.14 FT

ELEVATION OF GW PRIOR TO PURGING: 5.76 FT

THICKNESS OF LNAPL PRIOR TO PURGING : —.— FT

PID/Hnu READING IMMEDIATELY AFTER THE WELL CAP IS REMOVED: 0 PPM

pH: 5.45 TEMP: 17 C, SPECIFIC CONDUCTIVITY: 690 uS

DEPTH OF WELL: 16.52 FT D.O. - 3.7

HEIGHT OF WATER: 8.38 FT

EVACUATED GAL. H2O: 16.5 GAL (8.38 X .65 X 3 = 16.34)

PURGING START TIME: 9:23 END TIME: 9:41

PURGE METHOD: REDI-FLOW 2 INCH SUBMERSIBLE PUMP VARIABLE FLOW RATE OF <0.5 GPM TO >5.0 GPM

PURGE RATE (<0.5 GPM): 1 GPM

TOTAL VOLUME PURGED: 16.5 GAL.

DEPTH TO WATER AFTER PURGING AND BEFORE SAMPLING: 8.27 FT

DISSOLVED OXYGEN: 4.2 pH: 5.43 TEMP: 16 °C

SPECIFIC CONDUCTIVITY: 570 uS

SAMPLING METHOD: DEDICATED, DECONTAMINATED (IAW NJDEP FSPM 1992) TEFLON® BAILER

START TIME OF SAMPLING: 10:10 END TIME: 10:14

DISSOLVED OXYGEN: 3.6 pH: 5.30 TEMP: 14 °C

SPECIFIC CONDUCTIVITY: 540 uS

COMMENTS: _____

TOTAL ANALYTICAL SERVICES FOR A SAFE ENVIRONMENT

nytest environmental.

(516) 625-5500 FAX: (516) 625-1274

Chain of Custody Record

page # : 1 of 1

Client Name: Aguilar
Address: 30 Freeman Ave
Matawan, NJ 07747

Project Manager: D. Schmidt
Phone: 908-290-7500 FAX: 908-290-7506
Project Name: US Army, Ft Monmouth NJ GW 290
Project Number:
P.O. #:
Analytical Protocol: Deliverables:
Sampled By: D. Schmidt

Analysis Requested

Lab ID (Lab Use Only)	Sample ID (Maximum of 6 Characters)	Date Sampled	Time Sampled	Sample Location	No. of Containers	60X +15, Xylenes	+BTBE +TBA	Total Pb
100000	MW-1	11/8/94	1010	flushmount GW 290	3	2		1

Bin #'s In / Out (For Lab Use Only)

Login #:
Ship to:
Nytest Environmental Inc.
60 Seaview Blvd
Port Washington N.Y. 11050
Attn.: Sample Control
Date Shipped:
Carrier:
Air Bill #:
Cooler #:
C of C #:
SDG #:
NEI QT #:

Comments

Relinquished by:	Date / Time	Received by:	Date / Time
Sarah J Hallinan	11/9/94 0950	Jay Topfer	11/9 9:50
Print Name:		Print Name: Jay Topfer	
Relinquished by:	Date / Time	Received by:	Date / Time
Print Name:		Print Name:	
Relinquished by: Jay Topfer	Date / Time 11/9 5:30	Received by Laboratory:	Date / Time 11/9/94 17:30
Print Name: J Topfer		Print Name: PTRIPICIES	

Special Instructions :

CLIENT RETAINS YELLOW COPY ONLY

Lab Use Only

Custody Seals: Intact Broken Absent
Sample Rec'd in Good Condition? : Yes No
Sample Temperature: 4 Degrees Celsius
INSPECTED BY:
COMMENTS:

TOTAL ANALYTICAL SERVICES FOR A SAFE ENVIRONMENT

nytest environmental.

(516) 625-5500 FAX: (516) 625-1274

Chain of Custody Record

page # : 1 of 1

Client Name: Aguila

Address: 30 Freneau Ave
Matawan, NJ 07747

Project Manager: D. Schmidt

Phone: 908-290-7900 FAX: 908-290-7806

Project Name: US Army, Ft Monmouth, Charles Woods

Project Number:

P.O. #:

Analytical Protocol: _____ Deliverables: _____

Sampled By: D. Schmidt

Analysis Requested

Lab ID (Lab Use Only)	Sample ID (Maximum of 6 Characters)	Date Sampled	Time Sampled	Sample Location	No. of Containers	624 + 15 xylenes MIBK TBA	Total Lead	Bin #'s In / Out (For Lab Use Only)	Comments
	MW-3	11/5/94	1310	Shelter, 2562	3	2	1		
	MW-1	11/8/94	1413	Flush 2562	3	2	1		
	MW-2	11/8/94	1525	Flush, 2562	3	2	1		
	Trip B	11/8/94	1420			2			
200000									

Login #:

Ship to:
Nytest Environmental Inc.
60 Seaview Blvd
Port Washington N.Y. 11050
Attn.: Sample Control

Date Shipped:

Carrier:

Air Bill #:

Cooler #:

C of C #:

SDG #:

NEI QT #:

	Date / Time		Date / Time
Relinquished by: Sarah J Hulsland	11/9/94 0950	Received by: Jay Topfer	11/9 9:50
Print Name:		Print Name: Jay Topfer	
Relinquished by:		Received by:	
Print Name:		Print Name:	
Relinquished by: Jay Topfer	11/9 5:30	Received by Laboratory:	11/9/94 5:32
Print Name: J Topfer		Print Name:	

Special Instructions : _____

CLIENT RETAINS YELLOW COPY ONLY

Lab Use Only

Custody Seals: Intact Broken Absent

Sample Rec'd in Good Condition? Y N

Sample Temperature: _____ Degrees Celsius

INSPECTED BY:

COMMENTS:

nytest environmental.
(516) 625-5500 FAX: (516) 625-1274

94-1415

Chain of Custody Record

page # : 1 of 1

Client Name	Aguilar
Address	30 Frnezo Ave
	Matawan NJ 07747
Project Manager	D. Schmidt
Phone	908-290-7800 FAX 908 290-7806
Project Name	U.S. ARMY Ft MONMOUTH, NJ BLDG 108
Project Number	93-1-12-1934-29
P.O. #	
Analytical Protocol	Deliverables
Sampled By	D. Schmidt

Analysis Requested

Lab ID (Lab Use Only)	Sample ID (Maximum of 6 Characters)	Date Sampled	Time Sampled	Sample Location	No. of Containers	624+15+XYlenes +TBA+MBTE	Total Pb
	MW1	11/7/94	1049	Ft Monmouth	2		1
	MW2	11/7/94	1102	BLDG 108	2		1
	MW3	11/7/94	1040		2		1
	FIELD P	11/7/94	1030		2		1
	DUP	11/7/94	—		2		1
	TRIP BLD	11/7/94	—		2		1
9400000	MS MSD	11/7/94	1035		4		1

Bin #'s In / Out (For Lab Use Only)

Login #: 22899

Ship to:
Nytest Environmental Inc.
60 Seaview Blvd
Port Washington N.Y. 11050
Attn.: Sample Control

Date Shipped:
Carrier:
Air Bill #:
Cooler #:
C of C #:
SDG #:
NEI QT #:

Comments

	Date / Time		Date / Time
Relinquished by: Sarah J Hubbard	11/9/94 0950	Received by: Jay Topfer	11/9 9:50
Print Name:		Print Name: Jay Topfer	
Relinquished by:		Received by:	
Print Name:		Print Name:	
Relinquished by: Jay Topfer	11/9/94 5:30	Received by Laboratory: Pierre Pierre	11/9/94 5:30
Print Name: J Topfer		Print Name:	

Special Instructions :

CLIENT RETAINS YELLOW COPY ONLY

Lab Use Only

Custody Seals: Intact Broken Absent
Sample Bot'l in Good Condition? Y N
Sample Temperature: 4°C Preserved (Y/N)
INSPECTED BY:
COMMENTS:

U.S. ARMY FORT MONMOUTH

P.O. #: Ag. 21/94 0843

Chain of Custody

Project #: 93-11-30-1246-27
Customer:
Phone:

Sampler: D Schmidt / Ag. 12
Date / Time: 11/8/94 9:30

Site Name: Bld 290
NJDEP DICAR # 93-11-30-1246-27
MW Sampling

Analysis Parameters

Start:
Finish:

Preservation Method: HCL – VOA; HNO3 – Pb

Lab Sample ID Number	Date	Time	Customer Sample Location/ID Number	Sample Matrix	# of Bottles		VOA + TBA	Total Lead				Hnu	Remarks
1707.1	11/8/94	10/0	mw-1, 2930961	Aq	3	2		1				0	Sample temp 24°C
1702.4	11/7/94	1030	Field Blank	AQ	3	2		1				NB	
1702.5	11/7/94	NA	Duplicate	AQ	3	2		1				NA	
1705.4	11/8/94	NA	Trip Blank	AQ	3 2	2						NB	
1702.7	11/7/94	1035	ms/msd, mw-3	Aq	5	4		1				NB	

Hnu 510 861457
Hnu 0 ppm
Calibrated
8 ppm Benzene Surrogate
reads 8.6 ppm

Relinquished By (signature)	Date / Time	Received By (signature)	Shipped By:
[signature]	11-8-94 / 1630		

Relinquished By (signature)	Date / Time	Received for Lab by (signature):	Date / Time
[signature]	11/9/94 08:04	[signature]	11/9/94 / 0810

Note: A drawing depicting sample location should be attached or drawn on the reverse side of this chain of custody. NONE

SAI-ENV COC form 01 Page 1 of 1 Pages Rev. A Date: 02 Apr 93

Enviornmental Laboratory

Certification Number 13461

000004

[redacted] ENVIRONMENTAL INC.

INTERNAL CHAIN OF CUSTODY

Laboratory Person Breaking Field Seal on Sample Shuttle & Accepting Responsibility for Sample	NAME: Lien Hendes	TITLE: S.C.C.
Client: Aguilar Assoc.	Date Broken 11/09/94	Military Time Seal Broken: 1730
Login #: 22509	Analytical Parameter/Fraction: PG&E 624+15, Pb	

SAMPLE NO.	ALIQUOT/EXTRACT NO.	SAMPLE NO.	ALIQUOT/EXTRACT NO.
108-1	22509-01	750-4	22509-11
108-2	02	290-1	12
108-3	03	2562-3	13
108-FB	04	2562-1	14
108-DUP	05	2562-2	15
108-TB	06	2562-TB	16
108-3MS	07	296-6	17
108-3MSD	08	296-2 296-7	18 1c
750-1	09	296-3 296-8	2c 21
750-3	10	296-1	22

DATE	TIME	RELINQUISHED BY	RECIEVED BY	PURPOSE OF CHANGE OF CUST.
11/11/94	0800	PRINTED NAME M. Lam SIGNATURE M. Lam	PRINTED NAME SCarver SIGNATURE SCarver	PG&E 624+15
11/18/94	1500	PRINTED NAME M. LAM SIGNATURE M. Lam	PRINTED NAME H. Trujillo SIGNATURE Hall	Pb
11/30/94	1400	PRINTED NAME SCarver SIGNATURE SCarver	PRINTED NAME M. LAM SIGNATURE M. Lam	Storage
12/3/94	1000	PRINTED NAME H. Trujillo SIGNATURE Hall	PRINTED NAME M LAM SIGNATURE M. Lam	Storage
		PRINTED NAME SIGNATURE	PRINTED NAME SIGNATURE	
		PRINTED NAME SIGNATURE	PRINTED NAME SIGNATURE	
		PRINTED NAME SIGNATURE	PRINTED NAME SIGNATURE	
		PRINTED NAME SIGNATURE	PRINTED NAME SIGNATURE	

LABORATORY DELIVERABLES

		Check if Complete
1.	Cover page, Title page listing Lab Certification# facility name & address, & date of report	✓
2.	Table of Contents	✓
3.	Summary sheets listing analytical results for all targeted and non-targeted compounds	NA
4.	Summary Table cross-referencing field ID #'s vs. Lab ID #'s	✓
5.	Document bound, paginated and legible	✓
6.	Chain of Custody	✓
7.	Methodology Summary	✓
8.	Laboratory Chronicle and Holding Time check	✓
9.	Results submitted on a dry weight basis (if applicable)	NA
10.	Method Detection Limits	NA
11.	Lab certified by NJDEPE for parameters or appropriate category of parameters of the USEPA CLP	✓
12.	Non-Conformance Summary	✓

Laboratory Manager or Environmental Consultant's Signature

2/21/95
Date

GC/MS ANALYSIS CONFORMANCE/NON-CONFORMANCE SUMMARY FORMAT

		No	Yes
1.	Chromatograms Labeled/Compounds Identified (Field Samples and Method Blanks)	___	✓
2.	GC/MS Tune Specifications		
	a. BFB Meet Criteria	___	✓
	b. DFTPP Meet Criteria	___	NA
3.	GC/MS Tune Frequency - Performed every 24 hours for 600 series and 12 hours for 8000 series.	___	✓
4.	GC/MS Calibration - Initial Calibration performed within 30 days before sample analysis and continuing calibration performed within 24 hours of sample analysis for 600 series and 12 hours for 8000 series	___	✓
5.	GC/MS Calibration Requirements		
	a. Calibration Check Compounds	___	NA
	b. System Performance Check Compounds	___	NA

6. Blank Contamination - If yes, list compounds and concentrations in each blank:

   a. VOA fraction VBLKN04 - TBA - 2 ppb
   b. B/N Fraction _____
   c. Acid Fraction _____

7. Surrogate Recoveries Meet Criteria ___ ✓

   If not met, list those compounds and their recoveries which fall outside the acceptable range:

   a. VOA Fraction _____
   b. B/N Fraction _____
   c. Acid Fraction _____

   If not met, were the calculations checked and the results qualified as estimated? NA

8. Matrix Spike/Matrix Spike Duplicate Recoveries Meet Criteria (if not met, list these compounds and their recoveries which fall outside the acceptable range) ✓ ___

   a. VOA Fraction 1122-TCE - 162%
   b. B/N Fraction _____
   c. Acid Fraction _____

9. Internal Standard Area/Retention Time Shift Meet Criteria ___ NA ___

GC/MS ANALYSIS CONFORMANCE/NON-CONFORMANCE SUMMARY FORMAT (CONT.)

		No	Yes
10.	Extraction Holding Time Met	NA	
	If not met, list number of days exceeded each sample:		
11.	Analysis Holding Time Met		✓
	If not met, list number of days exceeded for each sample:		

Additional Comments: _____

Laboratory Manager: Yoni Bloy Date 2/21/95

METAL ANALYSIS CONFORMANCE/NON-CONFORMANCE SUMMARY FORMAT

		No	Yes
1.	Calibration Summary Meet Criteria	___	✓
2.	ICP Interference Check Sample Results Summary Submitted (if applicable) / Meet Criteria	___ NA	___
3.	Serial Dilution Summary Submitted (if applicable) / Meet Criteria	___ NA	___
4.	Laboratory Control Sample Summary Submitted (if applicable) / Meet Criteria	___	✓
5.	Blank Contamination - If yes, list compounds concentrations in each blank:	✓	___
6.	Matrix Spike/Matrix Spike Duplicate Recoveries Meet Criteria (If not met, list those compounds and their recoveries which fall outside the acceptable range)	___	✓
7.	Extraction Holding Time Met	___	✓
	If not met, list number of days exceeded for each sample:		
8.	Analysis Holding Time Met	___	✓
	If not met, list number of days exceeded for each sample:		

Additional Comments: _____

Laboratory Manager: Tom Berg Date: 2/21/95

Laboratory Chronicle

Client Name: Aguilar Associates — Log In No.: 22509E

Date(s) of Sample Collection: 11/07/94, 11/8/94

Date Received: 11/09/94

Sample ID: As per chain of custody

Organics Extraction:

1. Acids
2. Base/Neutrals
3. Pesticides/PCBs
4. Herbicides

Analysis: 11/14/94

1. Volatiles
2. Acids
3. Base/Neutrals
4. Pesticides/PCBs
5. Herbicides

Section Supervisor Review & Approval

Digestion - 11/21/94 Analysis - 12/09/94

Inorganics:

1. Metals
2. Cyanides
3. Phenols

Other Analysis:

Section Supervisor Review & Approval

Quality Control Supervisor Review & Approval

Dates are included for re-extractions and reanalysis.

000010

NARRATIVE DISCUSSION
VOLATILES - 22509E

Bldg. 290

INTRODUCTION

This narrative covers the analysis of three (3) samples in accordance with NEI Sop# 703 based on EPA Method 624.

HOLDING TIMES

The analytical holding time for this analysis was met.

CALIBRATIONS

All required minimum RRFs and maximum % RSD initial calibration requirements have been met in accordance with the Method.

QC CHECK SAMPLE

The % recoveries for four (4) out of thirty-one (31) compounds exceeded QC limits in the QC sample NO198.

Due to a laboratory oversight a second QC check was not analyzed.

METHOD BLANKS

The method blank associated with these samples met all method requirements.

SURROGATES

All surrogate recoveries met QC criteria.

MATRIX SPIKES

As requested, sample 108-3 was utilized in the MS/MSD series. The 1,1,2,2-Tetrachloroethane recovery in the MSD was above the advisory QC limits. The form 3 is included in this report.

INTERNAL STANDARDS

Internal Standard area responses/retention times summaries are not required.

SAMPLE COMMENTS

The TICs identified as "Unknown Siloxane" are most probably due to column degradation and not sample constituency.

NEI is reporting the results to our method detection limits (MDL's) rounded up to the nearest part per billion (ppb) in accordance with the guidance provided by NJDEP. These MDL's indicate that NEI did not detect any compounds above these levels.

Only one (1) field blank, 108FB, was collected with these samples. The field blank was collected on 11/07/94. The samples were collected on 11/08/94. The trip blank submitted with this report, 2562TB, was also collected on 11/08/94.

No further analytical problems were encountered.

NEI is reporting the results to our method detection limits (MDL's) rounded up to the nearest part per billion (ppb) in accordance with the guidance provided by NJDEP. These MDL's indicate that NEI did not detect any compounds above these levels.

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NARRATIVE DISCUSSION
INORGANICS- 22509E

All samples were analyzed as per required protocols.

The matrix spike recoveries and duplicate RPD values were within control limits.

NARRATIVE DISCUSSION
INORGANICS- 22509E

All samples were analyzed as per required protocols.

The matrix spike recoveries and duplicate RPD values were within control limits.

nytest environmental inc

I certify that this data package has been reviewed for the quality control and quality assurance measures for all analyzed methodologies.

Remo Gigante
Exec. Vice President

000016

METHODOLOGY SUMMARY

AQUEOUS METHODOLOGIES:	REF 1	REF 2	REF 3	REF 5
BNA, Pesticides/PCB's Extraction		3510/3520		
AA/ICP Sample Preparation	200.7			
Furnace Sample Preparation	200.0			
Mercury Sample Preparation	245.1			
Hexavalent Chromium Sample Preparation	218.5			
Clean-Up		3610/3620/3630/ 3640/3660		
Organochlorine Pesticide and PCB's by Gas Chromatography			608	505
Herbicides by Gas Chromatography			362	515.1
Purgeable Organics by GC/MS			624	524.2
Base/Neutral, Acids by GC/MS			625	525
2,3,7,8-TCDD by GC/MS			613/625	
BTEX			602	502.2
EDB/DBCP by Microextraction				504.1

NON-AQUEOUS METHODOLOGIES:	
BNA, Pesticides/PCB's Extraction	3550
AA/ICP Sample Preparation	3050
Furnace Sample Preparation	3020/3030/3050
Mercury Sample Preparation	7471
Clean-Up	3610/3620/3630/ 3640/3660

GC, Gas Chromatography/Mass Spectrometry:	
Purgeable Organics	8240/8021
Base/Neutral and Acid Extractables	8270
Organophosphorus Pesticides	8140
Organochlorine Pesticide and PCB's by Gas Chromatography	8080
BTEX	8020
Halogenated Purgeable Organics	8010

METHODOLOGY SUMMARY

INDUCTIVELY COUPLED PLASMA (ICP):	REFERENCE 1	REFERENCE 2
Aluminum	200.7	6010
Antimony	200.7	6010
Barium	200.7	6010
Beryllium	200.7	6010
Cadmium	200.7	6010
Calcium	200.7	6010
Chromium	200.7	6010
Cobalt	200.7	6010
Copper	200.7	6010
Iron	200.7	6010
Lead	200.7	6010
Magnesium	200.7	6010
Manganese	200.7	6010
Molybdenum	200.7	6010
Nickel	200.7	6010
Potassium	200.7	6010
Silver	200.7	6010
Sodium	200.7	6010
Tin	200.7	6010
Titanium	200.7	6010
Vanadium	200.7	6010
Zinc	200.7	6010
FURNACE AA:		
Antimony	204.1	7041
Arsenic	206.2	7060
Lead	239.2	7421
Selenium	270.2	7740
Thallium	279.2	7841
Tin	282.2	
Vanadium	286.2	7911
Mercury	245.1	7470/7471
ICAP:		
Priority Pollutants	200.7	6010/7060/ 7470/7740
TAL Metals	200.7	6010/7060/ 7470/7740
RCRA Metals	200.7	6010/7060/ 7470/7740

METHODOLOGY SUMMARY

REFERENCES:

(1) USEPA-600/4-79-020, Methods for Chemical Analysis of Water and Waste

(2) USEPA SW 846, Test Methods for Evaluating Solid Waste, Third Edition

(3) Federal Register 40 CFR Part 136, Vol.49, No.209 Test Parameters for the Analysis of Pollutants

(4) Federal Register Vol.51, No.216 Friday, 11/7/86, pp.40643-40652

(5) Method for the Determination of Organic Compounds in Drinking Water, EPA 500/4-88/039, Dec. 1988

(6) Standard Method for Examination of Water and Wastewater, 15 Edition 1980

Method Qualifiers for Organic Non-CLP Methodologies

Q Qualifier - Specified entries and their meanings as follows:

U - Indicates compound was analyzed for but was not detected. The sample quantitation limit is corrected for dilutions and for the moisture content for soil samples. If a sample extract can not be concentrated to the protocol - specific volume, this fact is also accounted for in reporting the sample quantitation limit. The number is the minimum detected limits for the sample.

J - Indicates an estimated volume. The flag is used either when estimating concentration for tentatively identified compounds where a 1:1 response is assumed, or when the mass spectral data indicates the presence of a compound that meets the identification criteria but the result is less than the sample quantitation limit but greater than zero.

N - Indicates presumptive evidence of a compound. This flag is used only for tentatively identified compounds, where the identification is based on a mass spectral library search. It is applied to all TIC results. For generic characterization of a TIC, such as chlorinated hydrocarbon, the **N** code is not used.

B - This flag is used when the analyte is found in the analyte is found in the associated blank as well as the sample. It indicates possible/probable blank contamination and warns the data user to take appropriate action. This flag is used for a TIC as well as for a positively identified target compound.

E - This flag identifies compounds whose concentrations exceeded the calibration range of the GC/MS instrument for that specific analysis.

D - This flag identifies all compounds identified in an analysis at a secondary dilution factor.

A - This flag indicates that a TIC is a suspected aldol condensation product.

Method Qualifiers for Inorganics

FORM I-IN includes fields for three types of results qualifiers. These qualifiers must be completed as follows:

* C (Concentration) qualifier -- Enter "B" if the reported value was obtained from a reading that was less than the Contract Required Detection Limit (CRDL) but greater than or equal to the Instrument Detection Limit (IDL). If the analyte was analyzed for but not detected, a "U" must be entered.

* Q Qualifier -- Specified entries and their meanings are as follows :

E - The reported value is estimated because of the presence of interference.

M - Duplicate precision not met (CV > 20%).

N - Spiked sample recovery not within control limits.

S - The reported value was determined by Method of Standard Addition (MSA).

W - Post-digestion spike for Furnace AA analysis is out of control limits (85-115%), while sample absorbance is less than 50% of spike absorbance.

* - Duplicate analysis not within control limits.

\+ - Correlation Coefficient for MSA is less than 0.995.

Entering "S", "W" or "+" is mutually exclusive.

* M (Method) qualifier - enter:

- "P" for ICP
- "A" for Flame AA
- "F" for Furnace AA
- "CV" for Cold Vapor .\A
- "AV" for Automated Cold Vapor AA
- "AS" for Semi-Automated Spectrophotometric
- "C" for Manual Spectrophotometric
- "T" for Titrimetric
- "NR" if the analyte is not required to be analyzed.

GC/MS Data

Volatile Data

1A

VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

108-FB

Lab Name: NYTEST ENV INC Contract: 9421415

Lab Code: NYTEST Case No.: 22509 SAS No.: SDG No.: ARMY1

Matrix: (soil/water) WATER Lab Sample ID: 2250904

Sample wt/vol: 5.0 (g/mL) ML Lab File ID: N0203.D

Level: (low/med) LOW Date Received: 11/09/94

% Moisture: not dec. _____ Date Analyzed: 11/14/94

Column: (pack/cap) CAP Dilution Factor: 1.0

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) UG/L	Q
74-87-3	Chloromethane	2	U
74-83-9	Bromomethane	1	U
75-01-4	Vinyl Chloride	1	U
75-00-3	Chloroethane	1	U
75-09-2	Methylene Chloride	4	
75-35-4	1,1-Dichloroethene	2	U
75-34-3	1,1-Dichloroethane	1	U
67-66-3	Chloroform	1	U
107-06-2	1,2-Dichloroethane	1	U
71-55-6	1,1,1-Trichloroethane	1	U
56-23-5	Carbon Tetrachloride	2	U
75-27-4	Bromodichloromethane	1	U
78-87-5	1,2-Dichloropropane	1	U
10061-01-5	cis-1,3-Dichloropropene	1	U
79-01-6	Trichloroethene	2	U
124-48-1	Dibromochloromethane	1	U
79-00-5	1,1,2-Trichloroethane	1	U
71-43-2	Benzene	1	U
10061-02-6	trans-1,3-Dichloropropene	1	U
75-25-2	Bromoform	1	U
127-18-4	Tetrachloroethene	3	U
79-34-5	1,1,2,2-Tetrachloroethane	2	U
108-88-3	Toluene	2	U
108-90-7	Chlorobenzene	2	U
100-41-4	Ethylbenzene	2	U
1330-20-7	Xylene (total)	6	U
75-69-4	Trichloromonofluoromethane	2	U
107-02-8	Acrolein	20	U
107-13-1	Acrylonitrile	2	U
75-65-0	Tertiary Butyl Alcohol	1	U
1634-34-4	Methyl Tertiary Butyl Ether	1	U
541-73-1	1,3-Dichlorobenzene	2	U
106-46-7	1,4-Dichlorobenzene	2	U
95-50-1	1,2-Dichlorobenzene	2	U

1A

VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

108-FB

Lab Name: NYTEST ENV INC Contract: 9421415

Lab Code: NYTEST Case No.: 22509 SAS No.: SDG No.: ARMY1

Matrix: (soil/water) WATER Lab Sample ID: 2250904

Sample wt/vol: 5.0 (g/mL) ML Lab File ID: N0203.D

Level: (low/med) LOW Date Received: 11/09/94

% Moisture: not dec. _____ Date Analyzed: 11/14/94

Column: (pack/cap) CAP Dilution Factor: 1.0

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) UG/L	Q
110-75-8	2-Chloroethylvinyl Ether	4	U
156-60-5	Trans, 1,2-Dichloroethene	1	U

1E

VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

108-FB

Lab Name: NYTEST ENV INC Contract: 9421415

Lab Code: NYTEST Case No.: 22509 SAS No.: SDG No.: ARMY1

Matrix: (soil/water) WATER Lab Sample ID: 2250904

Sample wt/vol: 5.0 (g/mL) ML Lab File ID: N0203.D

Level: (low/med) LOW Date Received: 11/09/94

% Moisture: not dec. _____ Date Analyzed: 11/14/94

Column: (pack/cap) CAP Dilution Factor: 1.0

Number TICs found: 1 CONCENTRATION UNITS: (ug/L or ug/Kg) UG/L

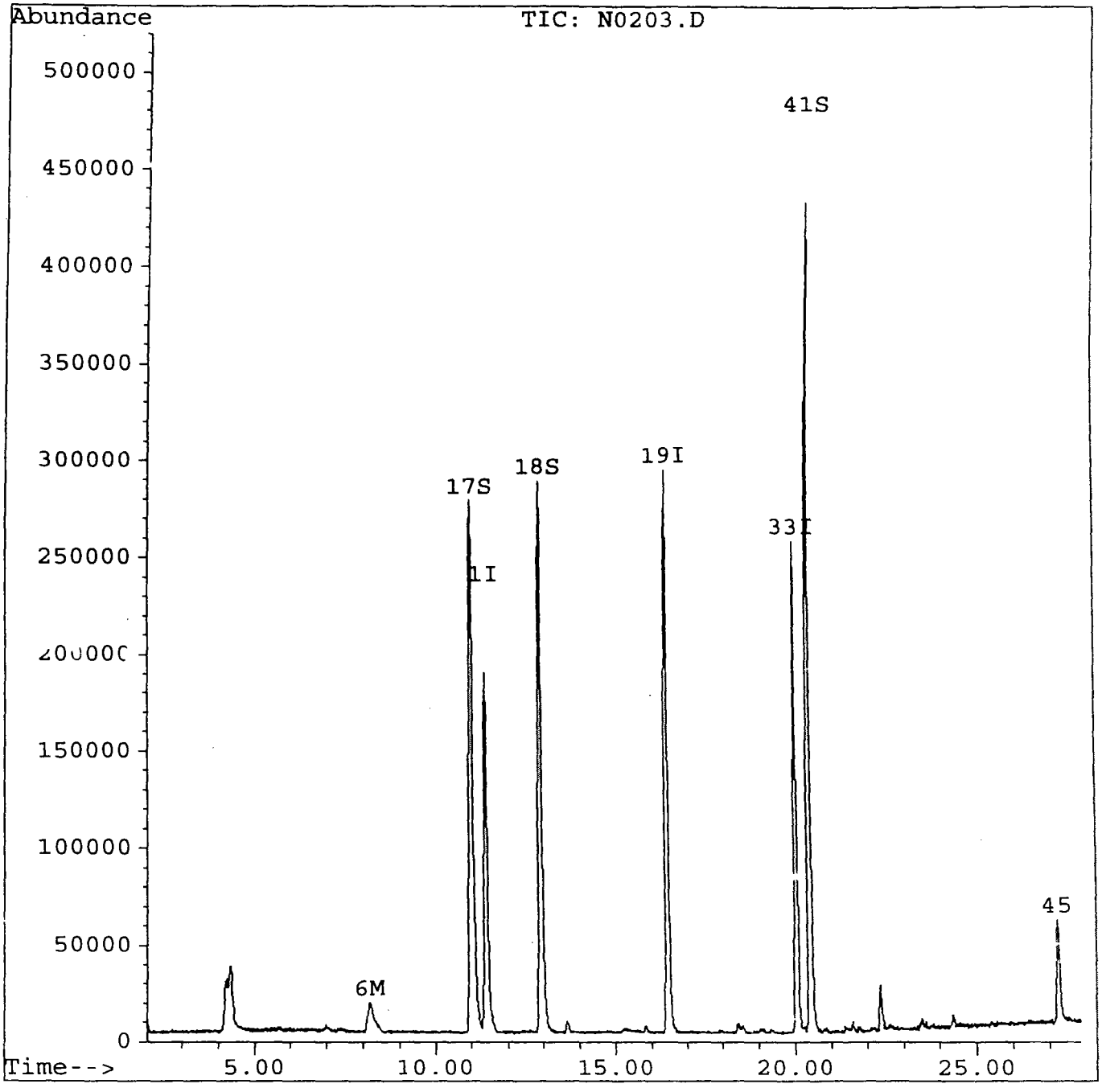
	CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1.		UNKNOWN HYDROCARBON	22.375	3	J
2.					
3.					
4.					
5.					
6.					
7.					
8.					
9.					
10.					
11.					
12.					
13.					
14.					
15.					
16.					
17.					
18.					
19.					
20.					
21.					
22.					
23.					
24.					
25.					
26.					
27.					
28.					
29.					
30.					

Quantitation Report

Data File : C:\HPCHEM\1\DATA\NOV1494\N0203.D
Acq Time : 14 Nov 94 13:44 pm
Sample : 2250904,108-FB,
Misc : 1,1,,,5,5,P624,R11-9-94
Quant Time: Dec 8 12:14 1994

Operator: LDS
Inst : HPN
Multiplr: 1.00

Method : c:\HPCHEM\1\METHODS\P624.M
Title : VOA Standards for 5 point calibration
Last Update : Tue Dec 06 02:26:13 1994
Response via : Multiple Level Calibration



000027

Quantitation Report

Data File : C:\HPCHEM\1\DATA\NOV1494\N0203.D
Acq Time : 14 Nov 94 13:44 pm
Sample : 2250904,108-FB,
Misc : 1,1,,,5,5,P624,R11-9-94
Quant Time: Dec 8 12:14 1994

Operator: LDS
Inst : HPN
Multiplr: 1.00

Method : c:\HPCHEM\1\METHODS\P624.M
Title : VOA Standards for 5 point calibration
Last Update : Tue Dec 06 02:26:13 1994
Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) CI01 Bromochloromethane	11.41	128	165759	30.00	ug/l	-0.07
19) 2-BROMO-1-CHLOROPROPANE	16.47	77	619401	30.00	ug/l	-0.04
33) 1,4-DICHLOROBUTANE	20.04	55	495243	30.00	ug/l	-0.03
System Monitoring Compounds						%Recovery
17) PENTAFLUOROBENZENE	11.02	168	781610	22.62	ug/l	75.40%
18) FLUOROBENZENE	12.95	96	917905	22.70	ug/l	75.67%
41) CS10 4-Bromofluorobenzene	20.40	95	527612	20.78	ug/l	69.28%
Target Compounds						Qvalue
6) C030 Methylene Chloride	8.19	84	51521	3.32	ug/l #	84
~~45) NAPHTHALENE~~	~~27.23~~	~~128~~	~~179646~~	~~3.31~~	~~ug/l~~	100

12-14

000028

(#) = qualifier out of range (m) = manual integration

1A

VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

290-1

Lab Name: NYTEST ENV INC Contract: 9421415

Lab Code: NYTEST Case No.: 22509 SAS No.: SDG No.: ARMY1

Matrix: (soil/water) WATER Lab Sample ID: 2250912

Sample wt/vol: 5.0 (g/mL) ML Lab File ID: N0211.D

Level: (low/med) LOW Date Received: 11/09/94

% Moisture: not dec. _____ Date Analyzed: 11/14/94

Column: (pack/cap) CAP Dilution Factor: 1.0

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) UG/L	Q
74-87-3	Chloromethane	2	U
74-83-9	Bromomethane	1	U
75-01-4	Vinyl Chloride	1	U
75-00-3	Chloroethane	1	U
75-09-2	Methylene Chloride	3	U
75-35-4	1,1-Dichloroethene	2	U
75-34-3	1,1-Dichloroethane	1	U
67-66-3	Chloroform	1	U
107-06-2	1,2-Dichloroethane	1	U
71-55-6	1,1,1-Trichloroethane	1	U
56-23-5	Carbon Tetrachloride	2	U
75-27-4	Bromodichloromethane	1	U
78-87-5	1,2-Dichloropropane	1	U
10061-01-5	cis-1,3-Dichloropropene	1	U
79-01-6	Trichloroethene	2	U
124-48-1	Dibromochloromethane	1	U
79-00-5	1,1,2-Trichloroethane	1	U
71-43-2	Benzene	1	U
10061-02-6	trans-1,3-Dichloropropene	1	U
75-25-2	Bromoform	1	U
127-18-4	Tetrachloroethene	3	U
79-34-5	1,1,2,2-Tetrachloroethane	2	U
108-88-3	Toluene	2	U
108-90-7	Chlorobenzene	2	U
100-41-4	Ethylbenzene	2	U
1330-20-7	Xylene (total)	6	U
75-69-4	Trichloromonofluoromethane	2	U
107-02-8	Acrolein	20	U
107-13-1	Acrylonitrile	2	U
75-65-0	Tertiary Butyl Alcohol	1	U
1634-34-4	Methyl Tertiary Butyl Ether	1	U
541-73-1	1,3-Dichlorobenzene	2	U
106-46-7	1,4-Dichlorobenzene	2	U
95-50-1	1,2-Dichlorobenzene	2	U

1A

VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

290-1

Lab Name: NYTEST ENV INC Contract: 9421415

Lab Code: NYTEST Case No.: 22509 SAS No.: SDG No.: ARMY1

Matrix: (soil/water) WATER Lab Sample ID: 2250912

Sample wt/vol: 5.0 (g/mL) ML Lab File ID: N0211.D

Level: (low/med) LOW Date Received: 11/09/94

% Moisture: not dec. _____ Date Analyzed: 11/14/94

Column: (pack/cap) CAP Dilution Factor: 1.0

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) UG/L	Q
110-75-8	2-Chloroethylvinyl Ether	4	U
156-60-5	Trans, 1,2-Dichloroethene	1	U

1E

VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

290-1

Lab Name: NYTEST ENV INC Contract: 9421415

Lab Code: NYTEST Case No.: 22509 SAS No.: SDG No.: ARMY1

Matrix: (soil/water) WATER Lab Sample ID: 2250912

Sample wt/vol: 5.0 (g/mL) ML Lab File ID: N0211.D

Level: (low/med) LOW Date Received: 11/09/94

% Moisture: not dec. _____ Date Analyzed: 11/14/94

Column: (pack/cap) CAP Dilution Factor: 1.0

Number TICs found: 3 CONCENTRATION UNITS: (ug/L or ug/Kg) UG/L

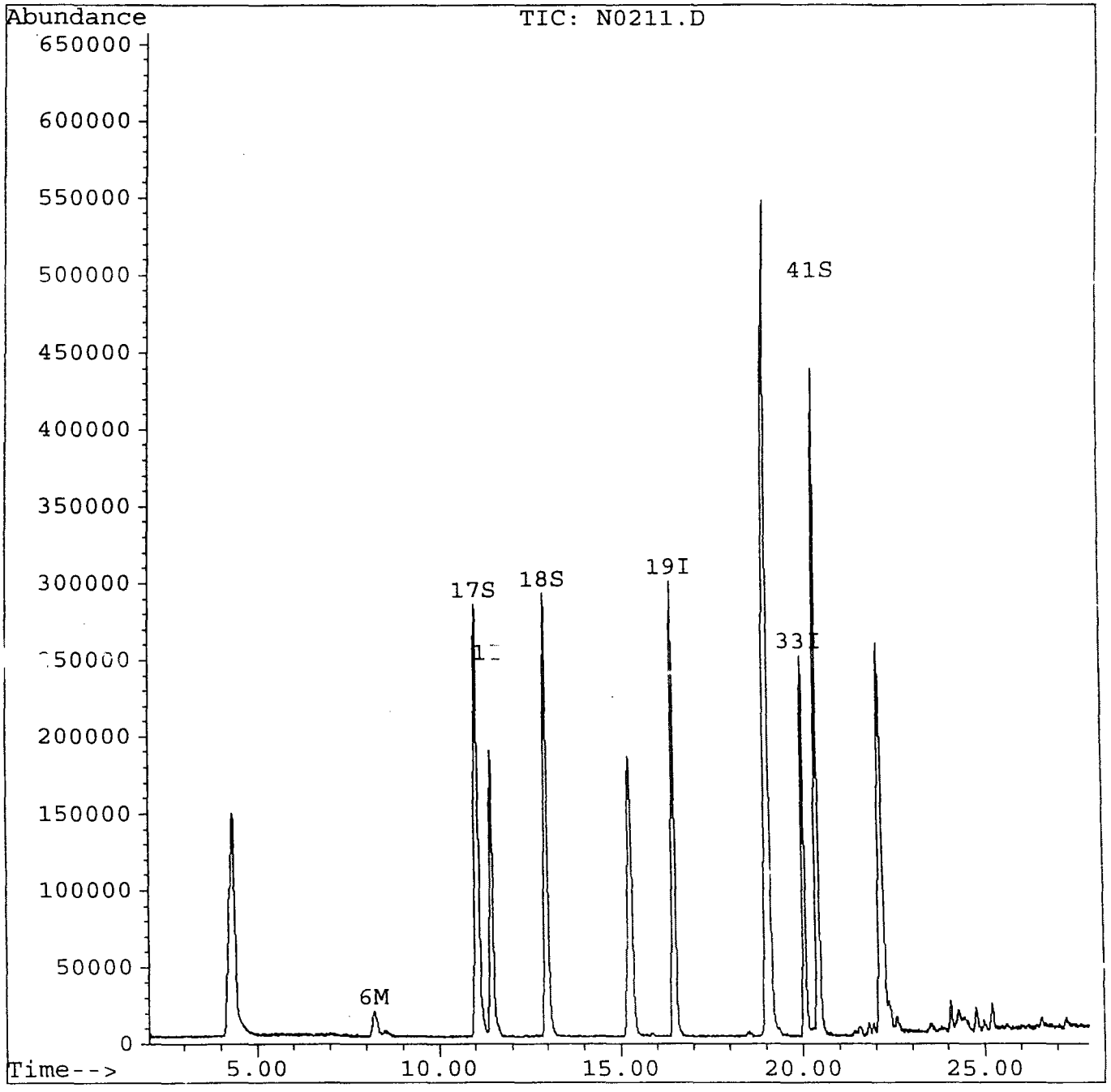
	CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1.		UNKNOWN SILOXANE	15.269	29	J
2.		UNKNOWN SILOXANE	19.069	110	J
3.		UNKNOWN SILOXANE	22.176	52	J
4.					
5.					
6.					
7.					
8.					
9.					
10.					
11.					
12.					
13.					
14.					
15.					
16.					
17.					
18.					
19.					
20.					
21.					
22.					
23.					
24.					
25.					
26.					
27.					
28.					
29.					
30.					

Quantitation Report

Data File : C:\HPCHEM\1\DATA\NOV1494\N0211.D
Acq Time : 14 Nov 94 18:11 pm
Sample : 2250912,290-1,
Misc : 1,1,,,5,5,P624,R11-9-94
Quant Time: Dec 8 12:34 1994

Operator: LDS
Inst : HPN
Multiplr: 1.00

Method : c:\HPCHEM\1\METHODS\P624.M
Title : VOA Standards for 5 point calibration
Last Update : Tue Dec 06 02:26:13 1994
Response via : Multiple Level Calibration



000032

Quantitation Report

Data File : C:\HPCHEM\1\DATA\NOV1494\N0211.D
Acq Time : 14 Nov 94 18:11 pm
Sample : 2250912,290-1,
Misc : 1,1,,,5,5,P624,R11-9-94
Quant Time: Dec 8 12:34 1994

Operator: LDS
Inst : HPN
Multiplr: 1.00

Method : c:\HPCHEM\1\METHODS\P624.M
Title : VOA Standards for 5 point calibration
Last Update : Tue Dec 06 02:26:13 1994
Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) CI01 Bromochloromethane	11.45	128	169672	30.00	ug/l	-0.03
19) 2-BROMO-1-CHLOROPROPANE	16.49	77	635282	30.00	ug/l	-0.01
33) 1,4-DICHLOROBUTANE	20.06	55	482239	30.00	ug/l	-0.01
System Monitoring Compounds						%Recovery
17) PENTAFLUOROBENZENE	11.05	168	789948	22.33	ug/l	74.45%
18) FLUOROBENZENE	12.97	96	943794	22.80	ug/l	76.01%
41) CS10 4-Bromofluorobenzene	20.43	95	532873	21.56	ug/l	71.86%
Target Compounds						Qvalue
6) C030 Methylene Chloride	8.21	84	20639	1.30	ug/l #	83

000033

(#) = qualifier out of range (m) = manual integration

1A

VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

2562-TB

Lab Name: NYTEST ENV INC Contract: 9421415

Lab Code: NYTEST Case No.: 22509 SAS No.: SDG No.: ARMY1

Matrix: (soil/water) WATER Lab Sample ID: 2250916

Sample wt/vol: 5.0 (g/mL) ML Lab File ID: N0215.D

Level: (low/med) LOW Date Received: 11/09/94

% Moisture: not dec. _____ Date Analyzed: 11/14/94

Column: (pack/cap) CAP Dilution Factor: 1.0

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) UG/L	Q
74-87-3	Chloromethane	2	U
74-83-9	Bromomethane	1	U
75-01-4	Vinyl Chloride	1	U
75-00-3	Chloroethane	1	U
75-09-2	Methylene Chloride	3	U
75-35-4	1,1-Dichloroethene	2	U
75-34-3	1,1-Dichloroethane	1	U
67-66-3	Chloroform	1	U
107-06-2	1,2-Dichloroethane	1	U
71-55-6	1,1,1-Trichloroethane	1	U
56-23-5	Carbon Tetrachloride	2	U
75-27-4	Bromodichloromethane	1	U
78-87-5	1,2-Dichloropropane	1	U
10061-01-5	cis-1,3-Dichloropropene	1	U
79-01-6	Trichloroethene	2	U
124-48-1	Dibromochloromethane	1	U
79-00-5	1,1,2-Trichloroethane	1	U
71-43-2	Benzene	1	U
10061-02-6	trans-1,3-Dichloropropene	1	U
75-25-2	Bromoform	1	U
127-18-4	Tetrachloroethene	3	U
79-34-5	1,1,2,2-Tetrachloroethane	2	U
108-88-3	Toluene	2	U
108-90-7	Chlorobenzene	2	U
100-41-4	Ethylbenzene	2	U
1330-20-7	Xylene (total)	6	U
75-69-4	Trichloromonofluoromethane	2	U
107-02-8	Acrolein	20	U
107-13-1	Acrylonitrile	2	U
75-65-0	Tertiary Butyl Alcohol	1	U
1634-34-4	Methyl Tertiary Butyl Ether	1	U
541-73-1	1,3-Dichlorobenzene	2	U
106-46-7	1,4-Dichlorobenzene	2	U
95-50-1	1,2-Dichlorobenzene	2	U

1A

EPA SAMPLE NO.

VOLATILE ORGANICS ANALYSIS DATA SHEET

2562-TB

Lab Name: NYTEST ENV INC Contract: 9421415

Lab Code: NYTEST Case No.: 22509 SAS No.: SDG No.: ARMY1

Matrix: (soil/water) WATER Lab Sample ID: 2250916

Sample wt/vol: 5.0 (g/mL) ML Lab File ID: N0215.D

Level: (low/med) LOW Date Received: 11/09/94

% Moisture: not dec. _____ Date Analyzed: 11/14/94

Column: (pack/cap) CAP Dilution Factor: 1.0

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) UG/L	Q
110-75-8	2-Chloroethylvinyl Ether	4	U
156-60-5	Trans, 1,2-Dichloroethene	1	U

1E

EPA SAMPLE NO.

VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

2562-TB

Lab Name: NYTEST ENV INC Contract: 9421415

Lab Code: NYTEST Case No.: 22509 SAS No.: SDG No.: ARMY1

Matrix: (soil/water) WATER Lab Sample ID: 2250916

Sample wt/vol: 5.0 (g/mL) ML Lab File ID: N0215.D

Level: (low/med) LOW Date Received: 11/09/94

% Moisture: not dec. _____ Date Analyzed: 11/14/94

Column: (pack/cap) CAP Dilution Factor: 1.0

Number TICs found: 0

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/L

	CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1.					
2.					
3.					
4.					
5.					
6.					
7.					
8.					
9.					
10.					
11.					
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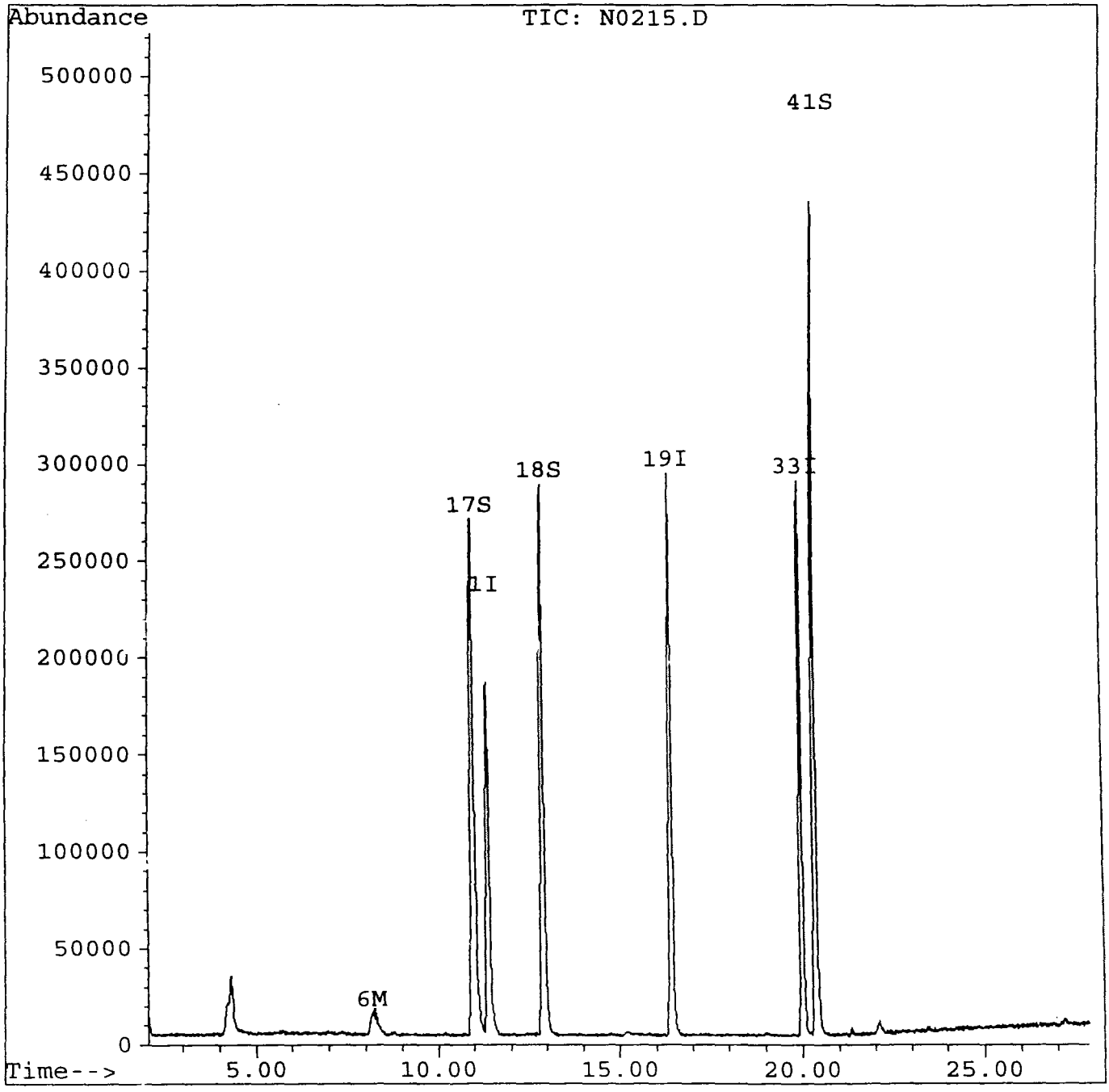
23.					
24.					
25.					
26.					
27.					
28.					
29.					
30.					

Quantitation Report

Data File : C:\HPCHEM\1\DATA\NOV1494\N0215.D
Acq Time : 14 Nov 94 20:26 pm
Sample : 2250916,2562-TB,
Misc : 1,1,,,5,5,P624,R11-9-94
Quant Time: Dec 8 12:40 1994

Operator: LDS
Inst : HPN
Multiplr: 1.00

Method : c:\HPCHEM\1\METHODS\P624.M
Title : VOA Standards for 5 point calibration
Last Update : Tue Dec 06 02:26:13 1994
Response via : Multiple Level Calibration



000037

Quantitation Report

Data File : C:\HPCHEM\1\DATA\NOV1494\N0215.D
Acq Time : 14 Nov 94 20:26 pm
Sample : 2250916,2562-TB,
Misc : 1,1,,,5,5,P624,R11-9-94
Quant Time: Dec 8 12:40 1994

Operator: LDS
Inst : HPN
Multiplr: 1.00

Method : c:\HPCHEM\1\METHODS\P624.M
Title : VOA Standards for 5 point calibration
Last Update : Tue Dec 06 02:26:13 1994
Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) CI01 Bromochloromethane	11.36	128	165852	30.00	ug/l	-0.12
19) 2-BROMO-1-CHLOROPROPANE	16.42	77	623345	30.00	ug/l	-0.09
33) 1,4-DICHLOROBUTANE	19.99	55	550392	30.00	ug/l	-0.08
System Monitoring Compounds						%Recovery
17) PENTAFLUOROBENZENE	10.97	168	767318	22.19	ug/l	73.98%
18) FLUOROBENZENE	12.90	96	921993	22.79	ug/l	75.97%
41) CS10 4-Bromofluorobenzene	20.36	95	514258	18.23	ug/l	60.76%
Target Compounds						Qvalue
6) C030 Methylene Chloride	8.17	84	21107	1.36	ug/l	87

000038

(#) = qualifier out of range (m) = manual integration

1A

VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

VBLKN04

Lab Name: NYTEST ENV INC Contract: 9421415

Lab Code: NYTEST Case No.: 22509 SAS No.: SDG No.: ARMY1

Matrix: (soil/water) WATER Lab Sample ID: VBLKN04

Sample wt/vol: 5.0 (g/mL) ML Lab File ID: N0199.D

Level: (low/med) LOW Date Received: 00/00/00

% Moisture: not dec. _____ Date Analyzed: 11/14/94

Column: (pack/cap) CAP Dilution Factor: 1.0

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) UG/L	Q
74-87-3	Chloromethane	2	U
74-83-9	Bromomethane	1	U
75-01-4	Vinyl Chloride	1	U
75-00-3	Chloroethane	1	U
75-09-2	Methylene Chloride	3	U
75-35-4	1,1-Dichloroethene	2	U
75-34-3	1,1-Dichloroethane	1	U
67-66-3	Chloroform	1	U
107-06-2	1,2-Dichloroethane	1	U
71-55-6	1,1,1-Trichloroethane	1	U
56-23-5	Carbon Tetrachloride	2	U
75-27-4	Bromodichloromethane	1	U
78-87-5	1,2-Dichloropropane	1	U
10061-01-5	cis-1,3-Dichloropropene	1	U
79-01-6	Trichloroethene	2	U
124-48-1	Dibromochloromethane	1	U
79-00-5	1,1,2-Trichloroethane	1	U
71-43-2	Benzene	1	U
10061-02-6	trans-1,3-Dichloropropene	1	U
75-25-2	Bromoform	1	U
127-18-4	Tetrachloroethene	3	U
79-34-5	1,1,2,2-Tetrachloroethane	2	U
108-88-3	Toluene	2	U
108-90-7	Chlorobenzene	2	U
100-41-4	Ethylbenzene	2	U
1330-20-7	Xylene (total)	6	U
75-69-4	Trichloromonofluoromethane	2	U
107-02-8	Acrolein	20	U
107-13-1	Acrylonitrile	2	U
75-65-0	Tertiary Butyl Alcohol	2	
1634-34-4	Methyl Tertiary Butyl Ether	1	U
541-73-1	1,3-Dichlorobenzene	2	U
106-46-7	1,4-Dichlorobenzene	2	U
95-50-1	1,2-Dichlorobenzene	2	U

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

VBLKN04

Lab Name: NYTEST ENV INC Contract: 9421415

Lab Code: NYTEST Case No.: 22509 SAS No.: SDG No.: ARMY1

Matrix: (soil/water) WATER Lab Sample ID: VBLKN04

Sample wt/vol: 5.0 (g/mL) ML Lab File ID: N0199.D

Level: (low/med) LOW Date Received: 00/00/00

% Moisture: not dec. _____ Date Analyzed: 11/14/94

Column: (pack/cap) CAP Dilution Factor: 1.0

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) UG/L	Q
110-75-8	2-Chloroethylvinyl Ether	4	U
156-60-5	Trans, 1,2-Dichloroethene	1	U

1E

VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

VBLKN04

Lab Name: NYTEST ENV INC Contract: 9421415

Lab Code: NYTEST Case No.: 22509 SAS No.: SDG No.: ARMY1

Matrix: (soil/water) WATER Lab Sample ID: VBLKN04

Sample wt/vol: 5.0 (g/mL) ML Lab File ID: N0199.D

Level: (low/med) LOW Date Received: 00/00/00

% Moisture: not dec. _____ Date Analyzed: 11/14/94

Column: (pack/cap) CAP Dilution Factor: 1.0

Number TICs found: 0

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/L

	CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1.					
2.					
3.					
4.					
5.					
6.					
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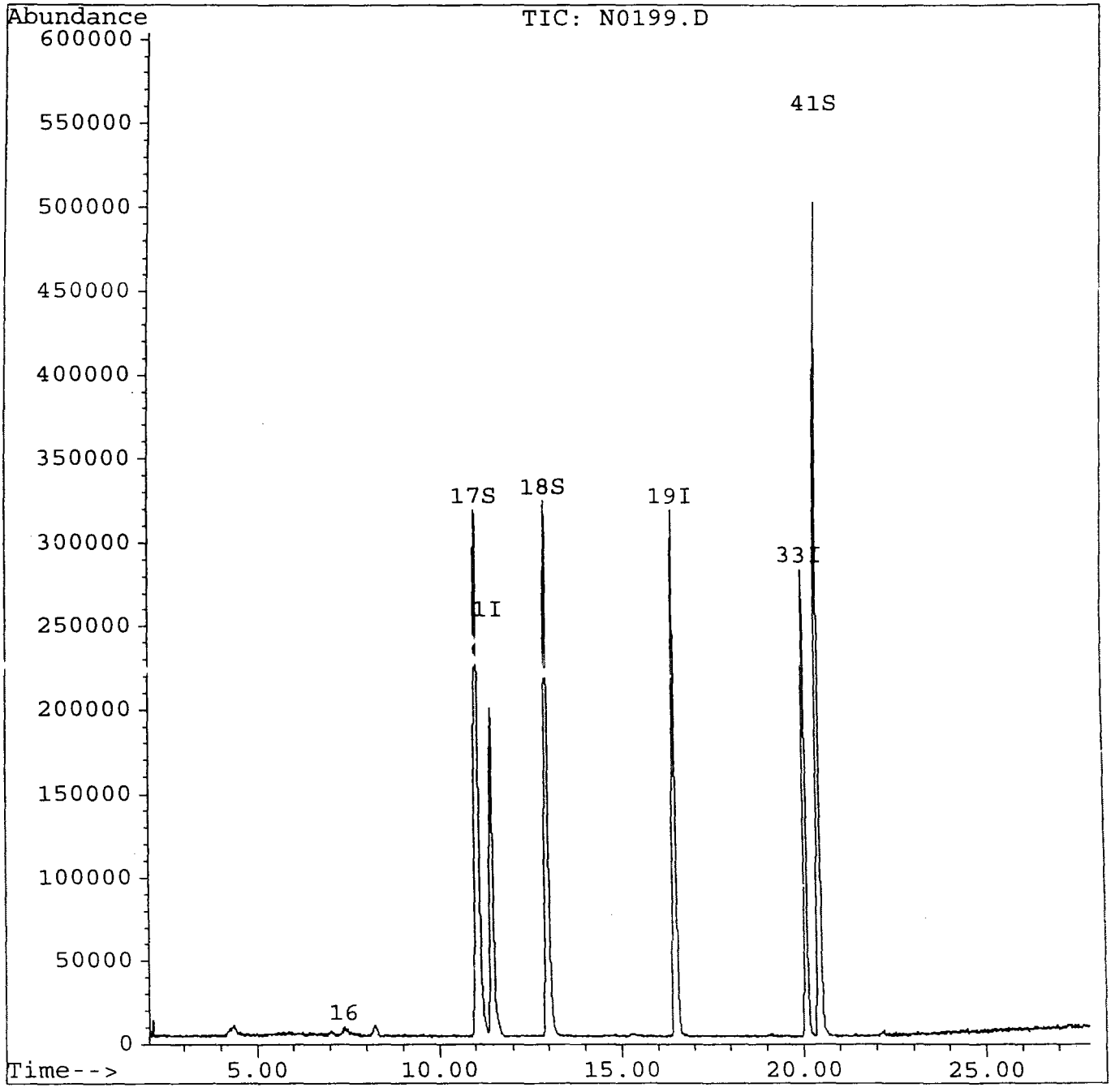
28.					
29.					
30.					

Quantitation Report

Data File : C:\HPCHEM\1\DATA\NOV1494\N0199.D
Acq Time : 14 Nov 94 11:31 am
Sample : VBLKN04,VBLKN04,
Misc : 1,,,,5,5,P624
Quant Time: Dec 8 12:08 1994

Operator: LDS
Inst : HPN
Multiplr: 1.00

Method : c:\HPCHEM\1\METHODS\P624.M
Title : VOA Standards for 5 point calibration
Last Update : Tue Dec 06 02:26:13 1994
Response via : Multiple Level Calibration



000042

Quantitation Report

Data File : C:\HPCHEM\1\DATA\NOV1494\N0199.D
Acq Time : 14 Nov 94 11:31 am
Sample : VBLKN04,VBLKN04,
Misc : 1,,,,5,5,P624
Quant Time: Dec 8 12:08 1994

Operator: LDS
Inst : HPN
Multiplr: 1.00

Method : c:\HPCHEM\1\METHODS\P624.M
Title : VOA Standards for 5 point calibration
Last Update : Tue Dec 06 02:26:13 1994
Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) CI01 Bromochloromethane	11.44	128	174658	30.00	ug/l	-0.04
19) 2-BROMO-1-CHLOROPROPANE	16.49	77	681811	30.00	ug/l	-0.01
33) 1,4-DICHLOROBUTANE	20.06	55	538079	30.00	ug/l	-0.02
System Monitoring Compounds						%Recovery
17) PENTAFLUOROBENZENE	11.04	168	890793	24.47	ug/l	81.55%
18) FLUCROBENZENE	12.97	96	1052247	24.70	ug/l	82.33%
41) CS10 4-Bromofluorobenzene	20.44	95	620849	22.51	ug/l	75.03%
Target Compounds						Qvalue
16) C177 TERTIARY BUTYL ALCOHO	7.38	59	29563	~~6.85~~	ug/l	100

2 ppb
us
12-14-94

000043

(#) = qualifier out of range (m) = manual integration

2A

WATER VOLATILE SURROGATE COMPOUND RECOVERY

Lab Name: NYTEST ENV INC Contract: 9421415

Lab Code: NYTEST Case No.: 22509 SAS No.: SDG No.: ARMY1

	EPA SAMPLE NO.	SUR1 (FB) #	SUR2 (BFB)#	SUR3 (PFB)#	OTHER	TOT OUT
01	QCCHECK	86	82	87		0
02	VBLKN04	93	85	92		0
03	108-1	88	84	86		0
04	108-2	88	79	87		0
05	108-3	87	79	86		0
06	108-FB	86	79	85		0
07	108-DUP	88	82	87		0
08	108-TB	87	80	86		0
09	108-3MS	87	76	85		0
10	108-3MSD	86	80	85		0
11	750-1	84	70	82		0
12	750-3	86	80	85		0
13	750-4	90	86	87		0
14	290-1	86	82	84		0
15	2562-3	86	75	86		0
16	2562-1	85	70	84		0
17	2562-2	85	69	84		0
18	2562-TB	86	69	84		0
19	296-6	84	68	83		0
20	296-2	85	69	82		0
21	296-3	87	78	86		0
22	296-8	93	71	87		0
23	296-1	84	74	81		0
24	296-7	87	80	84		0
25						

	QC LIMITS
SUR1 (FB) = Fluorobenzene	(70-140)
SUR2 (BFB) = Bromofluorobenzene	(60-150)
SUR3 (PFB) = Pentafluorobenzene	(70-140)

Column to be used to flag recovery values

* Values outside of contract required QC limits

D Surrogates diluted out

nytest

RECOVERY REPORT

Client Name:
Sample Matrix: LIQUID
Lab Smp Id: 2250907
Level: LOW
Data Type: MS DATA
SpikeList File: QCMS.spk
Method File: /chem/HPN.i/22509.b/624.m
Misc Info:

Client SDG: ARMY1
Fraction: VOA
Client Smp ID: 108-3MS
Operator: LDS
SampleType: MS
Quant Type: ISTD

SPIKE	COMPOUND	CONC ADDED ug/l	CONC RECOVERED ug/l	% RECOVERED	LIMITS
1	Chloromethane	10	8	84.79	0-273
3	Bromomethane	10	12	121.47	0-242
4	Vinyl Chloride	10	9	92.95	0-251
5	Chloroethane	10	12	119.98	14-230
6	Methylene Chloride	10	12	121.12	0-221
9	1,1-Dichloroethene	10	12	120.19	0-234
11	1,1-Dichloroethane	10	13	127.58	59-155
13	Trans, 1,2-Dichlor	10	12	123.47	54-156
14	Chloroform	10	13	131.21	51-138
15	1,2-Dichloroethane	10	13	133.47	49-155
18	Trichloromonofluor	10	11	106.23	17-181
23	1,1,1-Trichloroeth	10	12	117.25	52-162
24	Carbon Tetrachlori	10	11	113.33	70-140
26	Bromodichlorometha	10	12	121.27	35-155
27	1,2-Dichloropropan	10	12	121.82	0-210
28	cis-1,3-Dichloropr	10	13	131.08	0-227
29	Trichloroethene	10	11	112.34	71-157
30	Dibromochlorometha	10	12	119.42	53-149
31	1,1,2-Trichloroeth	10	13	130.49	52-150
33	Benzene	10	12	123.04	37-151
34	trans-1,3-Dichloro	10	12	125.82	17-183
36	Bromoform	10	13	127.32	45-169
37	2-Chloroethylvinyl	10	12	121.83	0-305
40	Tetrachloroethene	10	10	102.44	64-148
42	1,1,2,2-Tetrachlor	10	16	156.75	46-157
44	Toluene	10	13	128.56	47-150
45	Chlorobenzene	10	11	107.79	37-160
46	Ethylbenzene	10	9	88.28	37-162
52	1,3-Dichlorobenzen	10	11	106.99	59-156
53	1,4-Dichlorobenzen	10	10	96.14	18-190
54	1,2-Dichlorobenzen	10	13	132.56	18-190

nytest

RECOVERY REPORT

Client Name:
Sample Matrix: LIQUID
Lab Smp Id: 2250907
Level: LOW
Data Type: MS DATA
SpikeList File: QCMS.spk
Method File: /chem/HPN.i/22509.b/624.m
Misc Info:

Client SDG: ARMY1
Fraction: VOA
Client Smp ID: 108-3MS
Operator: LDS
SampleType: MS
Quant Type: ISTD

SURROGATE COMPOUND	CONC ADDED ug/l	CONC RECOVERED ug/l	% RECOVERED	LIMITS
$ 32 Pentafluorobenzene	30	25	84.67	70-1..
$ 43 Fluorobenzene	30	26	87.14	70-140
$ 47 Bromofluorobenzene	30	23	75.82	60-150

Data File: /chem/HPN.i/22509.b/N0207.d
Report Date: 15-Dec-1994 13:02

nytest

RECOVERY REPORT

Client Name:
Sample Matrix: LIQUID
Lab Smp Id: 2250908
Level: LOW
Data Type: MS DATA
SpikeList File: QCMS.spk
Method File: /chem/HPN.i/22509.b/624.m
Misc Info:

Client SDG: ARMY1
Fraction: VOA
Client Smp ID: 108-3MSD
Operator: LDS
SampleType: MSD
Quant Type: ISTD

SPIKE	COMPOUND	CONC ADDED ug/l	CONC RECOVERED ug/l	% RECOVERED	LIMITS
1	Chloromethane	10	10	95.98	0-2
3	Bromomethane	10	12	122.21	0-242
4	Vinyl Chloride	10	10	96.32	0-251
5	Chloroethane	10	12	120.96	14-230
6	Methylene Chloride	10	13	128.86	0-221
9	1,1-Dichloroethene	10	12	122.93	0-234
11	1,1-Dichloroethane	10	13	126.36	59-155
13	Trans, 1,2-Dichlor	10	12	124.98	54-156
14	Chloroform	10	13	133.38	51-138
15	1,2-Dichloroethane	10	13	132.51	49-155
18	Trichloromonofluor	10	11	113.32	17-181
23	1,1,1-Trichloroeth	10	13	126.26	52-162
24	Carbon Tetrachlori	10	12	118.85	70-140
26	Bromodichlorometha	10	12	125.76	35-155
27	1,2-Dichloropropan	10	13	129.27	0-210
28	cis-1,3-Dichloropr	10	13	127.86	0-227
29	Trichloroethene	10	12	117.50	71-157
30	Dibromochlorometha	10	12	123.75	53-149
31	1,1,2-Trichloroeth	10	14	135.49	52-150
33	Benzene	10	13	127.54	37-151
34	trans-1,3-Dichloro	10	13	126.89	17-183
36	Bromoform	10	12	125.03	45-169
37	2-Chloroethylvinyl	10	13	129.28	0-305
40	Tetrachloroethene	10	11	110.89	64-148
42	1,1,2,2-Tetrachlor	10	16	161.70*	46-157
44	Toluene	10	13	127.12	47-150
45	Chlorobenzene	10	12	118.34	37-160
46	Ethylbenzene	10	11	106.44	37-162
52	1,3-Dichlorobenzen	10	12	119.84	59-156
53	1,4-Dichlorobenzen	10	10	103.60	18-190
54	1,2-Dichlorobenzen	10	14	142.02	18-190

nytest

RECOVERY REPORT

Client Name:
Sample Matrix: LIQUID
Lab Smp Id: 2250908
Level: LOW
Data Type: MS DATA
SpikeList File: QCMS.spk
Method File: /chem/HPN.i/22509.b/624.m
Misc Info:

Client SDG: ARMY1
Fraction: VOA
Client Smp ID: 108-3MSD
Operator: LDS
SampleType: MSD
Quant Type: ISTD

SURROGATE COMPOUND	CONC ADDED ug/l	CONC RECOVERED ug/l	% RECOVERED	LIMITS
$ 32 Pentafluorobenzene	30	25	84.88	70-140
$ 43 Fluorobenzene	30	26	85.98	70-140
$ 47 Bromofluorobenzene	30	24	79.61	60-150

4A

VOLATILE METHOD BLANK SUMMARY

EPA SAMPLE NO.

VBLKN04

Lab Name: NYTEST ENV INC Contract: 9421415

Lab Code: NYTEST Case No.: 22509 SAS No.: SDG No.: ARMY1

Lab File ID: N0199.D Lab Sample ID: VBLKN04

Date Analyzed: 11/14/94 Time Analyzed: 1131

Matrix: (soil/water) WATER Level: (low/med) LOW

Instrument ID: HPN

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES, MS AND MSD

	EPA SAMPLE NO.	LAB SAMPLE ID	LAB FILE ID	TIME ANALYZED
01	QCCHECK	QCCHECK	N0198.D	1058
02	108-1	2250901	N0200.D	1205
03	108-2	2250902	N0201.D	1238
04	108-3	2250903	N0202.D	1311
05	108-FB	2250904	N0203.D	1344
06	108-DUP	2250905	N0204.D	1418
07	108-TB	2250906	N0205.D	1451
08	108-3MS	2250907	N0206.D	1524
09	108-3MSD	2250908	N0207.D	1558
10	750-1	2250909	N0208.D	1631
11	750-3	2250910	N0209.D	1704
12	750-4	2250911	N0210.D	1738
13	290-1	2250912	N0211.D	1811
14	2562-3	2250913	N0212.D	1845
15	2562-1	2250914	N0213.D	1919
16	2562-2	2250915	N0214.D	1953
17	2562-TB	2250916	N0215.D	2026
18	296-6	2250917	N0216.D	2100
19	296-2	2250918	N0217.D	2135
20	296-3	2250920	N0219.D	2234
21	296-8	2250921	N0220.D	2307
22	296-1	2250922	N0221.D	2340
23	296-7	2250919	N0222.D	0013
24				
25				
26				
27				
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29				
30				

COMMENTS:

5A

VOLATILE ORGANIC INSTRUMENT PERFORMANCE CHECK
BROMOFLUOROBENZENE (BFB)

Lab Name: NYTEST ENV INC Contract: 9421415

Lab Code: NYTEST Case No.: 22509 SAS No.: SDG No.: ARMY1

Lab File ID: N0163.D BFB Injection Date: 11/08/94

Instrument ID: HPN BFB Injection Time: 1152

Matrix:(soil/water) WATER Level:(low/med) LOW Column:(pack/cap) CAP

m/e	ION ABUNDANCE CRITERIA	% RELATIVE ABUNDANCE
50	15.0 - 40.0% of mass 95	15.4
75	30.0 - 60.0% of mass 95	38.4
95	Base peak, 100% relative abundance	100.0
96	5.0 - 9.0% of mass 95	7.3
173	Less than 2.0% of mass 174	0.0 (0.0)1
174	Greater than 50.0% of mass 95	67.3
175	5.0 - 9.0% of mass 174	5.2 (7.7)1
176	Greater than 95.0%, but less than 101.0% of mass 174	65.8 (97.8)1
177	5.0 - 9.0% of mass 176	4.5 (6.8)2

1-Value is % mass 174 2-Value is % mass 176

THIS CHECK APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS, AND STANDARDS:

	EPA SAMPLE NO.	LAB SAMPLE ID	LAB FILE ID	DATE ANALYZED	TIME ANALYZED
01	VSTD005	VSTD005	N0165.D	11/08/94	1237
02	VSTD010	VSTD010	N0166.D	11/08/94	1310
03	VSTD030	VSTD030	N0167.D	11/08/94	1343
04	VSTD050	VSTD050	N0168.D	11/08/94	1416
05	VSTD200	VSTD200	N0169.D	11/08/94	1450
06					
07					
08					
09					
10					
11					
12					
13					
14					
15					
16					
17					
18					
19					
20					
21					
22					

5A
VOLATILE ORGANIC INSTRUMENT PERFORMANCE CHECK
BROMOFLUOROBENZENE (BFB)

Lab Name: NYTEST ENV INC Contract: 9421415

Lab Code: NYTEST Case No.: 22509 SAS No.: SDG No.: ARMY1

Lab File ID: N0197.D BFB Injection Date: 11/14/94

Instrument ID: HPN BFB Injection Time: 1028

Matrix:(soil/water) WATER Level:(low/med) LOW Column:(pack/cap) CAP

m/e	ION ABUNDANCE CRITERIA	% RELATIVE ABUNDANCE
50	15.0 - 40.0% of mass 95	15.1
75	30.0 - 60.0% of mass 95	40.3
95	Base peak, 100% relative abundance	100.0
96	5.0 - 9.0% of mass 95	7.0
173	Less than 2.0% of mass 174	0.0 (0.0)1
174	Greater than 50.0% of mass 95	69.8
175	5.0 - 9.0% of mass 174	5.6 (8.0)1
176	Greater than 95.0%, but less than 101.0% of mass 174	69.8 (100.0)1
177	5.0 - 9.0% of mass 176	4.0 (5.8)2

1-Value is % mass 174 2-Value is % mass 176

THIS CHECK APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS, AND STANDARDS:

	EPA SAMPLE NO.	LAB SAMPLE ID	LAB FILE ID	DATE ANALYZED	TIME ANALYZED
01	QCCHECK	QCCHECK	N0198.D	11/14/94	1058
02	VBLKN04	VBLKN04	N0199.D	11/14/94	1131
03	108-1	2250901	N0200.D	11/14/94	1205
04	108-2	2250902	N0201.D	11/14/94	1238
05	108-3	2250903	N0202.D	11/14/94	1311
06	108-FB	2250904	N0203.D	11/14/94	1344
07	108-DUP	2250905	N0204.D	11/14/94	1418
08	108-TB	2250906	N0205.D	11/14/94	1451
09	108-3MS	2250907	N0206.D	11/14/94	1524
10	108-3MSD	2250908	N0207.D	11/14/94	1558
11	750-1	2250909	N0208.D	11/14/94	1631
12	750-3	2250910	N0209.D	11/14/94	1704
13	750-4	2250911	N0210.D	11/14/94	1738
14	290-1	2250912	N0211.D	11/14/94	1811
15	2562-3	2250913	N0212.D	11/14/94	1845
16	2562-1	2250914	N0213.D	11/14/94	1919
17	2562-2	2250915	N0214.D	11/14/94	1953
18	2562-TB	2250916	N0215.D	11/14/94	2026
19	296-6	2250917	N0216.D	11/14/94	2100
20	296-2	2250918	N0217.D	11/14/94	2135
21	296-3	2250920	N0219.D	11/14/94	2234
22	296-8	2250921	N0220.D	11/14/94	2307

5A
VOLATILE ORGANIC INSTRUMENT PERFORMANCE CHECK
BROMOFLUOROBENZENE (BFB)

Lab Name: NYTEST ENV INC Contract: 9421415

Lab Code: NYTEST Case No.: 22509 SAS No.: SDG No.: ARMY1

Lab File ID: N0197.D BFB Injection Date: 11/14/94

Instrument ID: HPN BFB Injection Time: 1028

Matrix:(soil/water) WATER Level:(low/med) LOW Column:(pack/cap) CAP

m/e	ION ABUNDANCE CRITERIA	% RELATIVE ABUNDANCE
50	15.0 - 40.0% of mass 95	15.1
75	30.0 - 60.0% of mass 95	40.3
95	Base peak, 100% relative abundance	100.0
96	5.0 - 9.0% of mass 95	7.0
173	Less than 2.0% of mass 174	0.0 (0.0)1
174	Greater than 50.0% of mass 95	69.8
175	5.0 - 9.0% of mass 174	5.6 (8.0)1
176	Greater than 95.0%, but less than 101.0% of mass 174	69.8 (100.0)1
177	5.0 - 9.0% of mass 176	4.0 (5.8)2

1-Value is % mass 174 2-Value is % mass 176

THIS CHECK APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS, AND STANDARDS:

	EPA SAMPLE NO.	LAB SAMPLE ID	LAB FILE ID	DATE ANALYZED	TIME ANALYZED
01	296-1	2250922	N0221.D	11/14/94	2340
02	296-7	2250919	N0222.D	11/15/94	0013
03					
04					
05					
06					
07					
08					
09					
10					
11					
12					
13					
14					
15					
16					
17					
18					
19					
20					
21					
22					

6A

VOLATILE ORGANICS INITIAL CALIBRATION DATA

Lab Name: NYTEST ENV INC Contract: 9421415

Lab Code: NYTEST Case No.: 22509 SAS No.: SDG No.: ARMY1

Instrument ID: HPN Calibration Date(s): 11/08/94

Matrix:(soil/water) WATER Level:(low/med) LOW Column:(pack/cap) CAP

Max $\overline{\%RSD}$ for CCC(*) = 35.0%

LAB FILE ID: RRF005 =N0165.D RRF010 =N0166.D
RRF030=N0167.D RRF050=N0168.D RRF200=N0169.D

COMPOUND	RRF005	RRF010	RRF030	RRF050	RRF200	$\overline{RRF}$	% RSD
Chloromethane	1.569	1.346	1.370	1.315	1.288	1.378	8.1
Bromomethane	1.967	1.807	1.716	1.642	1.596	1.746	8.4
Vinyl_Chloride	1.965	1.785	1.822	1.747	1.789	1.821	4.6
Chloroethane	1.133	1.109	1.001	1.021	0.975	1.048	6.6
Methylene_Chloride	3.913	2.873	2.154	2.030	1.854	2.565	33.0
1,1-Dichloroethene	2.105	1.877	1.918	1.889	1.855	1.929	5.2
1,1-Dichloroethane	3.845	3.464	3.518	3.485	3.589	3.580	4.3
Chloroform	4.126	3.689	3.562	3.506	3.588	3.694	6.8
1,2-Dichloroethane	2.024	1.859	1.799	1.741	1.835	1.852	5.7
1,1,1-Trichloroethane	0.939	0.857	0.859	0.827	0.866	0.869	4.8
Carbon_Tetrachloride	0.814	0.736	0.760	0.751	0.778	0.768	3.9
Bromodichloromethane	1.081	0.977	0.971	0.953	1.010	0.998	5.1
1,2-Dichloropropane	0.722	0.658	0.658	0.647	0.686	0.674	4.5
cis-1,3-Dichloropropene	0.983	0.873	0.869	0.846	0.897	0.893	5.9
Trichloroethene	0.712	0.643	0.653	0.652	0.669	0.666	4.1
Dibromochloromethane	0.883	0.816	0.812	0.803	0.843	0.831	3.9
1,1,2-Trichloroethane	0.629	0.570	0.564	0.548	0.578	0.578	5.3
Benzene	1.788	1.593	1.612	1.576	1.647	1.643	5.2
trans-1,3-Dichloropropene	0.792	0.710	0.708	0.688	0.726	0.725	5.5
Bromoform	0.709	0.646	0.644	0.645	0.565	0.642	7.9
Tetrachloroethene	0.921	0.805	0.826	0.835	0.902	0.858	5.9
1,1,2,2-Tetrachloroethane	1.209	1.036	1.013	0.986	0.000	1.061	9.5
Toluene	2.017	1.801	1.774	1.746	1.820	1.832	5.8
Chlorobenzene	2.021	1.789	1.801	1.797	1.970	1.876	5.9
Ethylbenzene	0.976	0.823	0.799	0.835	0.905	0.868	8.3
Xylene_(total)	1.192	1.065	1.058	1.066	1.152	1.106	5.5
Trichloromonofluoromethane	3.860	3.429	3.389	3.310	3.404	3.479	6.3
Acrolein	0.139	0.100	0.096	0.087	0.072	0.099	25.4
Acrylonitrile	0.792	0.744	0.714	0.764	0.695	0.742	5.2
Tertiary_Butyl_Alcohol	1.994	1.821	3.284	3.221	1.915	2.447	30.2
Methyl_Tertiary_Butyl_Ether	4.512	4.110	3.702	3.671	3.551	3.909	10.2
1,3-Dichlorobenzene	1.776	1.527	1.544	1.574	1.335	1.551	10.1
1,4-Dichlorobenzene	1.808	1.663	1.553	1.722	1.257	1.600	13.3
1,2-Dichlorobenzene	1.663	1.436	1.449	1.469	0.000	1.504	7.1
2-Chloroethylvinyl_Ether	0.722	0.658	0.658	0.647	0.686	0.674	4.5
Trans,_1,2-Dichloroethene	2.270	1.921	1.880	1.838	1.838	1.950	9.4
Fluorobenzene	7.222	6.388	6.218	6.099	6.397	6.465	6.8

000053

6A

VOLATILE ORGANICS INITIAL CALIBRATION DATA

Lab Name: NYTEST ENV INC Contract: 9421415

Lab Code: NYTEST Case No.: 22509 SAS No.: SDG No.: ARMY1

Instrument ID: HPN Calibration Date(s): 11/08/94

Matrix:(soil/water) WATER Level:(low/med) LOW Column:(pack/cap) CAP

Max $\overline{\%RSD}$ for CCC(*) = 35.0%

LAB FILE ID: RRF005 =N0165.D RRF010 =N0166.D
RRF030=N0167.D RRF050=N0168.D RRF200=N0169.D

COMPOUND	RRF005	RRF010	RRF030	RRF050	RRF200	$\overline{RRF}$	% RSD
Bromofluorobenzene	1.548	1.324	1.291	1.240	1.371	1.355	8.7
Pentafluorobenzene	6.013	5.464	5.291	5.277	5.534	5.516	5.4

060054

nytest

RECOVERY REPORT

Client Name:
Sample Matrix: LIQUID
Lab Smp Id: QCCHECK
Level: LOW
Data Type: MS DATA
SpikeList File: QCCHK.spk
Method File: /chem/HPN.i/22509.b/624.m
Misc Info:

Client SDG: ARMY1
Fraction: VOA
Client Smp ID: QCCHECK
Operator: LDS
SampleType: MS
Quant Type: ISTD

SPIKE	COMPOUND	CONC ADDED ug/l	CONC RECOVERED ug/l	% RECOVERED	LIMITS
1	Chloromethane	10	9	93.65	0-2[illegible]
3	Bromomethane	10	12	120.55	14-186
4	Vinyl Chloride	10	9	93.76	4-196
5	Chloroethane	10	12	117.23	38-162
6	Methylene Chloride	10	13	128.27	60-140
9	1,1-Dichloroethene	10	12	115.97	50-150
11	1,1-Dichloroethane	10	12	121.88	72-128
13	Trans, 1,2-Dichlor	10	12	121.73	70-130
14	Chloroform	10	13	129.64	68-132
15	1,2-Dichloroethane	10	13	129.33	68-132
18	Trichloromonofluor	10	11	108.43	48-152
23	1,1,1-Trichloroeth	10	12	123.83	75-125
24	Carbon Tetrachlori	10	12	120.99	73-127
26	Bromodichlorometha	10	12	123.55	66-134
27	1,2-Dichloropropan	10	12	123.82	34-166
28	cis-1,3-Dichloropr	10	13	128.31	24-176
29	Trichloroethene	10	12	119.42	66-134
30	Dibromochlorometha	10	12	124.96	68-132
31	1,1,2-Trichloroeth	10	12	124.50	71-129
33	Benzene	10	12	123.90	64-136
34	trans-1,3-Dichloro	10	13	128.05	50-150
36	Bromoform	10	13	130.91*	71-129
37	2-Chloroethylvinyl	10	12	123.83	0-224
40	Tetrachloroethene	10	11	114.93	74-126
42	1,1,2,2-Tetrachlor	10	16	159.06*	60-140
44	Toluene	10	13	126.54*	74-126
45	Chlorobenzene	10	12	118.24	66-134
46	Ethylbenzene	10	12	117.79	59-141
52	1,3-Dichlorobenzen	10	12	124.53	73-127
53	1,4-Dichlorobenzen	10	11	108.55	63-137
54	1,2-Dichlorobenzen	10	15	148.69*	63-127

nytest

RECOVERY REPORT

Client Name:
Sample Matrix: LIQUID
Lab Smp Id: QCCHECK
Level: LOW
Data Type: MS DATA
SpikeList File: QCCHK.spk
Method File: /chem/HPN.i/22509.b/624.m
Misc Info:

Client SDG: ARMY1
Fraction: VOA
Client Smp ID: QCCHECK
Operator: LDS
SampleType: MS
Quant Type: ISTD

SURROGATE COMPOUND	CONC ADDED ug/l	CONC RECOVERED ug/l	% RECOVERED	LIMITS
$ 32 Pentafluorobenzene	30	26	86.62	70-140
$ 43 Fluorobenzene	30	26	85.75	70-140
$ 47 Bromofluorobenzene	30	25	82.02	60-150

Metals Data

NYTEST ENVIRONMENTAL INC

INORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

108-FB

Lab Name: NYTEST_ENV_INC._____ Contract: 9421415___

Lab Code: NYTEST Login No.: 22509_ QC Report No.22509_

Matrix (soil/water): WATER Lab Sample ID: 250904____
Level (low/high) : LOW Date Received: 11/09/94
Percent Solids : __0.0

Concentration Units (ug/L or mg/kg dry weight): UG/L_

CAS No.	Analyte	Concentration	C	Q	M
7439-92-1	Lead	4.1	B		F

CODES :
P: ICP; F : GFAA; CV: Cold Vapor; AS: Automated Spectrophotometric
Note: A "U" in the "C" (Concentration) column indicates the analyte was not detected in this sample; "B" = Sample value greater than Instrument Detection Limit, but less than reporting limit; "NR" = Not Required.

Comments:
108-FB_____

000058

NYTEST ENVIRONMENTAL INC.

INORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

290-1

Lab Name: NYTEST_ENV_INC._____ Contract: 9421415___

Lab Code: NYTEST Login No.: 22509_ QC Report No.22509_

Matrix (soil/water): WATER
Level (low/high) : LOW
Percent Solids : __0.0

Lab Sample ID: 250912____
Date Received: 11/09/94

Concentration Units (ug/L or mg/kg dry weight): UG/L_

CAS No.	Analyte	Concentration	C	Q	M
7439-92-1	Lead	17.0			F

CODES :
P: ICP; F : GFAA; CV: Cold Vapor; AS: Automated Spectrophotometric
Note: A "U" in the "C" (Concentration) column indicates the analyte was not detected in this sample; "B" = Sample value greater than Instrument Detection Limit, but less than reporting limit; "NR" = Not Required.

Comments:
290-1

NYTEST ENVIRONMENTAL INC.

INITIAL AND CONTINUING CALIBRATION VERIFICATION

Lab Name: NYTEST_ENV_INC._____ Contract: 9421415___

Lab Code: NYTEST Login No.: 22509_ QC Report No.: 22509_

Initial Calibration Source: SPEX_____

Continuing Calibration Source: SPEX_____

Concentration Units: ug/L

Analyte	Initial Calibration True	Found	%R(1)	Continuing Calibration True	Found	%R(1)	Found	%R(1)	M
Lead	90.0	87.00	96.7	50.0	49.70	99.4	49.70	99.4	F

NR : Analyte Not Required

NYTEST ENVIRONMENTAL INC.

INITIAL AND CONTINUING CALIBRATION VERIFICATION

Lab Name: NYTEST_ENV_INC._____ Contract: 9421415___

Lab Code: NYTEST Login No.: 22509_ QC Report No.: 22509_

Initial Calibration Source: SPEX_____

Continuing Calibration Source: SPEX_____

Concentration Units: ug/L

Analyte	Initial Calibration True	Found	%R(1)	Continuing Calibration True	Found	%R(1)	Found	%R(1)	M
Lead				50.0	51.10	102.2	49.50	99.0	F

NR : Analyte Not Required

NYTEST ENVIRONMENTAL INC.

CRDL STANDARD FOR AA AND ICP

Lab Name: NYTEST_ENV_INC._____ Contract: 9421415___

Lab Code: NYTEST Login No.: 22509_ QC Report No. : 22509__

AA CRDL Standard Source: SPEX_____

ICP CRDL Standard Source: _____

Concentration Units: ug/L

	CRDL Standard for AA			CRDL Standard for ICP				
					Initial			Final
Analyte	True	Found	%R	True	Found	%R	Found	%R
Lead	3.0	2.60	86.7					

NEI FORM 3 - (1/94)

000062

ILMO2.1

NYTEST ENVIRONMENTAL INC.

ANALYTICAL AND METHOD BLANK SUMMARY

Lab Name: NYTEST_ENV_INC._____ Contract: 9421415___

Lab Code: NYTEST Login No.: 22509_ QC Report No.: 22509__

Preparation Blank Matrix (soil/water): WATER

Preparation Blank Concentration Units (ug/L or mg/kg): UG/L_

Analyte	Initial Calib. Blank (ug/L)	C	Continuing Calibration Blank (ug/L) 1	C	2	C	3	C	Prepa-ration Blank	C	M
Lead	3.0	U	3.0	U	3.0	U	3.0	U	3.000	U	F

NR = Analyte Not Requested

NYTEST ENVIRONMENTAL INC.

ANALYTICAL AND METHOD BLANK SUMMARY

Lab Name: NYTEST_ENV_INC._____ Contract: 9421415___

Lab Code: NYTEST Login No.: 22509_ QC Report No.: 22509__

Preparation Blank Matrix (soil/water): _____

Preparation Blank Concentration Units (ug/L or mg/kg): _____

Analyte	Initial Calib. Blank (ug/L)	C	Continuing Calibration Blank (ug/L) 1	C	2	C	3	C	Prepa-ration Blank	C	M
Lead			3.0	U							F

NR = Analyte Not Requested

NYTEST ENVIRONMENTAL INC.

MATRIX SPIKE RECOVERY DATA SHEET

SAMPLE NO.

108-3MSD

Lab Name: NYTEST_ENV_INC. Contract: 9421415

Lab Code: NYTEST Login No.: 22509 QC Report No. : 22509

Matrix (soil/water): WATER Level (low/med): LOW

% Solids for Sample: 0.0

Concentration Units (ug/L or mg/kg dry weight): UG/L

Analyte	Control Limit %R	Spiked Sample Result (SSR)	C	Sample Result (SR)	C	Spike Added (SA)	%R	Q	M
Lead	75-125	35.4000		16.7000		20.00	93.5		F

Comments:

108-3MSD

NR : Analyte Not Required

NEI FORM 7 - (1/94)

000065

NYTEST ENVIRONMENTAL INC.

DUPLICATES

SAMPLE NO.

108-3MS

Lab Name: NYTEST_ENV_INC. Contract: 9421415

Lab Code: NYTEST Login No.: 22509 QC Report No. : 22509

Matrix (soil/water): WATER Level (low/med): LOW

% Solids for Sample: 0.0 % Solids for Duplicate: 0.0

Concentration Units (ug/L or mg/kg dry weight): UG/L

Analyte	Control Limit	Sample (S)	C	Duplicate (D)	C	RPD	Q	M
Lead	5.0	16.7000		16.5000		1.2		F

NR : Analyte Not Requested

NEI FORM 9 (1/94)

NYTEST ENVIRONMENTAL INC.

LABORATORY CONTROL SAMPLE

Lab Name: NYTEST_ENV_INC._____ Contract: 9421415___

Lab Code: NYTEST Login No.: 22509_ QC Report No.: 22509_

Solid LCS Source: _____

Aqueous LCS Source: SPEX_____

Analyte	Aqueous (ug/L) True	Found	%R	Solid (mg/kg) True	Found	C	Limits		%R
Lead	90.0	78.60	87.3						

NYTEST ENVIRONMENTAL INC.

Instrument Detection Limits (Quarterly)

Lab Name: NYTEST_ENV_INC._____ Contract: 9421415___

Lab Code: NYTEST Login No.: 22509_ QC Report No.: 22509__

ICP ID Number: _____ Date in Effect: 10/01/94

Flame AA ID Number : _____

Furnace AA ID Number : 3030#1_____

Analyte	Wave-length (nm)	Back-ground	CRDL (ug/L)	IDL (ug/L)	M
Lead	283.30	BZ	5	3.0	F

Comments:

_____

_____

_____

_____

NEI FORM 13 - (1/94)

000068

Project No.: 9421415
Log in No. : 22633B
P.O. No. : Pending
Date : 2/21/95
SDG No. : Army 3
NJDEPE Case #:93-11-30-1246-27

ANALYTICAL DATA REPORT
PACKAGE FOR

Aguilar Associates

30 Freneau Avenue

Matawan, NJ 07747

ATTN: Darryl Schmitt
REF: US Army Fort Monmouth, Well# and NJDEPE Reg# 1-2930961
Sample Location Bldg. 290

LABORATORY NUMBER	SAMPLE IDENTIFICATION	TYPE OF SAMPLE

S E E N E X T P A G E

WE CERTIFY THAT THIS REPORT IS A TRUE REPORT OF RESULTS OBTAINED FROM OUR TESTS OF THIS MATERIAL.

RESPECTFULLY SUBMITTED,
NYTEST ENVIRONMENTAL INC.

NYS Lab ID. #10195
NJ Cert. #73469

REMO GIGANTE
EXEC. VICE PRESIDENT

NYTEST ENVIRONMENTAL Inc

LABORATORY NUMBER	SAMPLE IDENTIFICATION	WELL #	TYPE OF SAMPLE
2263312	290-1	1-2930961	Water
2263307	296-TB	-	Water
2263308	296-FB	-	Water

U.S. Army, Fort Monmouth
Sample ID Cross Reference Table

Building #	NJDEPE MW#	Sample ID
290	1-2930961	290-1

Table of Contents

TOTAL ANALYTICAL SERVICES FOR A SAFE ENVIRONMENT

NEI nytest environmental.

(516) 625-5500 FAX: (516) 625-1274

Chain of Custody Record

Client Name: Aguilar Assoc. & Cons., Inc.
Address: 30 Freneau Ave.
Matawan, NJ 07747

Project Manager: D. Schmitt
Phone: (908) 290-7800 FAX: (908) 290-7806
Project Name: Fort Monmouth Bldg. 290
Project Number: 93-11-30-1246-27
P.O. #:
Analytical Protocol: Deliverables:
Sampled By: D. Schmitt

Login #:
Ship to:
Nytest Environmental Inc.
60 Seaview Blvd
Port Washington N.Y. 11050
Attn.: Sample Control
Date Shipped:
Carrier:
Air Bill #:
Cooler #:
C of C #:
SDG #:
NEI QT #:

Analysis Requested

Lab ID (Lab Use Only)	Sample ID (Maximum of 6 Characters)	Date Sampled	Time Sampled	Sample Location	No. of Containers	GC/MS, no hits, TBA, MTBE	Total Lead								Comments
	MW-1	11/29/94	13:05	Bldg 290	3	✓	✓								

Bin #'s In / Out (For Lab Use Only)

000001

Relinquished by	Date / Time	Received by	Date / Time
Darryl W. Schmitt (signature); Print Name: Darryl W. Schmitt	11/30/94 1200	(signature); Print Name: T. Rymanowski	11/30/94 16:00
Relinquished by: ; Print Name:	Date / Time	Received by: ; Print Name:	Date / Time
(signature); Print Name: T. Rymanowski	11/30/94 18:20	Received by Laboratory: (signature); Print Name: Pieri Pierides	11/30/94 1800

Lab Use Only
Custody Seals: Intact Broken Absent
Sample Rec'd In Good Condition? : Y N
Sample Temperature: 6 Degrees Celcius
INSPECTED BY: PP
COMMENTS:

Special Instructions :

CLIENT RETAINS YELLOW COPY ONLY

BLDG.#: 290 W#: 1 NJDEPE WELL ID # 2930961

U.S. ARMY FORT MONMOUTH

MONITORING WELL SAMPLING DATASHEET

DATE: 11-29-94

IJO#94-0843 A

SAMPLING CONTRACTOR: Aguilar Associates Inc.

LABORATORY: NYTEST Environmental Inc. CERT #:73469

SAMPLERS NAMES: D. Schmitt, N. Vrasek, S. Parizzi

WEATHER CONDITIONS: Mid 40's, cold, windy

ELEVATION OF CASING SURVEY MARK: 13.90

TOTAL DEPTH OF WELL FROM TOP OF SURVEYORS MARK: 16.47 FT

DEPTH FROM SURVEYORS MARK TO SCREEN: 2.0 FT

LENGTH OF SCREENED SECTION: 10.5 FT.

DEPTH TO WATER PRIOR TO PURGING AND SAMPLING: 7.15 FT

ELEVATION OF GW PRIOR TO PURGING: 6.75 FT

THICKNESS OF LNAPL PRIOR TO PURGING : —.— FT

PID/Hnu READING IMMEDIATELY AFTER THE WELL CAP IS REMOVED: 0 PPM

pH: 5.74 TEMP: 13 C, SPECIFIC CONDUCTIVITY: 670 µS D.O.- 3.2

DEPTH OF WELL: 16.47 FT

HEIGHT OF WATER: 9.32 FT

EVACUATED GAL. H2O: 18 GAL (9.32 X .65 X 3 = 18.17)

PURGING START TIME: 11:20 END TIME: 11:42

PURGE METHOD: REDI-FLOW 2 INCH SUBMERSIBLE PUMP VARIABLE FLOW RATE OF <0.5 GPM TO >5.0 GPM

PURGE RATE (<0.5 GPM): 1 GPM

TOTAL VOLUME PURGED: 18 GAL.

DEPTH TO WATER AFTER PURGING AND BEFORE SAMPLING: 7.19 FT

DISSOLVED OXYGEN: 3.4 pH: 5.50 TEMP: 16 °C

SPECIFIC CONDUCTIVITY: 650 µS

SAMPLING METHOD: DEDICATED, DECONTAMINATED (IAW NJDEP FSPM 1992) TEFLON® BAILER

START TIME OF SAMPLING: 12:00 END TIME: 12:05

DISSOLVED OXYGEN: 2.8 pH: 5.61 TEMP: 15 °C

SPECIFIC CONDUCTIVITY: 620 µS

COMMENTS: _____

U.S. ARMY FORT MONMOUTH

P.O. #: Aguifer / 94-0843

Chain of Custody

Project #: 93-11-3D-1246-27	Sampler: D. Schmitt - Aguifer Associates	Date: 11/29/94	Time: 1100
Customer: C. Appleby SELFM-PW-EV	Site Name: Bldg. 270		
Phone: (908) 532-6224	DECAP #: 93-11-30-1246-27		
	MW Samplings		

Analysis Parameters

Start:

Finish:

Preservation Method

Lab Sample ID Number	Date/Time		Customer Sample Location/ID Number	Sample Matrix	# of Bottles	BTEX, MTBE, Total Lead	Hnu (ppm)	Remarks
1742.1	11/29/94	12:05	MW-1, 293096I	AQ	3	✓ ✓	0	Sample Kept <4°C
1741.7	11/29/94	9:53	Field Blank	AQ	3	✓ ✓		
1741.8	11/29/94	9:35	Trip Blank	AQ	2	✓		
1741.9	11/29/94	-	Duplicate	AQ	3	✓ ✓		
1741.10	11/29/94	10:11	mw-6, ms/ds, 293097S	AQ	3	✓ ✓		
								Hnu SN 801497
								Hnu = 0 ppm
								Calibrated 11/23/94
								5ppm Benzene Standard reads 5.0 ppm span 9.5

D. S.

Relinquished By (signature)	Date / Time	Received By (signature)	Shipped By:
Daryl W. Schmitt	11/30/94 1100	Sarah Hubbard	
Relinquished By (signature)	Date / Time	Received for Lab by (signature):	Date / Time

Note: A drawing depicting sample location should be attached or drawn on the reverse side of this chain of custody. NONE

Enviornmental Laboratory

Certification Number: 1246

TOTAL ANALYTICAL SERVICES FOR A SAFE ENVIRONMENT

nytest environmental.

(516) 625-5500 FAX: (516) 625-1274

12/1 44214

Chain of Custody Record

page # : 1 of 1

Client Name	Aguilar Assoc. & Cons., Inc.
Address	30 Frenchan Ave.
	Matawan, NJ 07747
Project Manager	D. Schmitt
Phone	(908) 290-7800 FAX (908) 290-7806
Project Name	Fort Monmouth Bldg 296
Project Number	93-11-2-1200-13
P.O. #	
Analytical Protocol	Deliverables
Sampled By	D. Schmitt

Analysis Requested

Lab ID (Lab Use Only)	Sample ID (Maximum of 6 Characters)	Date Sampled	Time Sampled	Sample Location	No. of Containers	624 VOAs + Xylenes, MTBE, TBA	Total Lead	Bin #'s In / Out (Lab Use Only)	Comments
1	mw-1	11/30/94	11:50	Stickup Bld 296	3	✓	✓		
2	mw-2		11:40	Street Bld 296					
3	mw-3		10:00	Stickup Bld 296					
4	mw-6		10:11	Stickup Bld 296					
5	mw-7		10:30	fence Bld 296					
6	mw-8		9:20	Street Bld 296					
7	Field B		9:53	Field BLANK			✓		
8	Trip B		9:25	Trip blank					
9	Dup 1-C		—	Duplicate			✓		
10,11	mw-6 ms		10:11	MS/MSD			✓		

Login #: 22637

Ship to:
Nytest Environmental Inc.
60 Seaview Blvd
Port Washington N.Y. 11050
Attn.: Sample Control

Date Shipped:
Carrier:
Air Bill #:
Cooler #:
C of C #:
SDG #:
NEI QT #:

	Date / Time		Date / Time
Relinquished by: Darryl W. Schmitt	11/30/94 1600	Received by: T. Rymanowski	11/30/94 16:00
Print Name: Darryl W. Schmitt		Print Name:	
Relinquished by:		Received by:	
Print Name:		Print Name:	
Relinquished by: T. Rymanowski	11/30/94 18:20	Received by Laboratory: Pieri Pierides	11/30/94 1820
Print Name:		Print Name:	

Lab Use Only

Custody Seals: Intact / Broken / Absent
Sample Rec'd in Good Condition? Y / N
Sample Temperature: 6 Degrees Celsius
INSPECTED BY:
COMMENTS:

Special Instructions :

CLIENT RETAINS YELLOW COPY ONLY

NYTEST ENVIRONMENTAL INC.

INTERNAL CHAIN OF CUSTODY

Laboratory Person Breaking Field Seal on Sample Shuttle & Accepting Responsibility for Sample	NAME: Pierr Perdes	TITLE: SCO
Client: Aguilar Associates	Date Broken 11/30/94	Military Time Seal Broken: 1320
Login #: 22633	Analytical Parameter/Fraction: P624+15, Pb, 8080PEST, 8150	

SAMPLE NO.	ALIQUOT/EXTRACT NO.	SAMPLE NO.	ALIQUOT/EXTRACT NO.
296-1	22633-01	296-MSD	22633-11
296-2	02	290-1	12
296-3	03	108-1	13
296-6	04	108-2	14
296-7	05	108-3	15
296-8	06	750-1	16
296-FB	07	750-3	17
296-TB	08	750-4	18
296-DUP	09	2562-1	19
296-MS	10	2562-2	20
2502-3	21		
2044-1	22		
2044-2	23		
2044-3	24		
2044-FB	25		
2044-DUP	26		
2044-TB	27		

DATE	TIME	RELINQUISHED BY	RECIEVED BY	PURPOSE OF CHANGE OF CUST
12/1/94	1152	PRINTED NAME M. LANI SIGNATURE M. Lan	PRINTED NAME P. BCALES SIGNATURE	P624+15
12/1/94	1310	PRINTED NAME M. LANI SIGNATURE M. Lan	PRINTED NAME C VOSS SIGNATURE C Voss	8080PEST, 8150
12/6/94	1500	PRINTED NAME C VOSS SIGNATURE C Voss	PRINTED NAME P. PERALES SIGNATURE	STORAGE
12/6/94	1443	PRINTED NAME M LANI SIGNATURE M. Lan	PRINTED NAME A. Frimpong SIGNATURE A. Frimpong	Pb
12/8/94	1500	PRINTED NAME A. Frimpong SIGNATURE A. Frimpong	PRINTED NAME M. LANI SIGNATURE M. Lan	STORAGE
12/13/94	0800	PRINTED NAME P. PERALES SIGNATURE	PRINTED NAME M. LANI SIGNATURE M. Lan	STORAGE
		PRINTED NAME SIGNATURE	PRINTED NAME SIGNATURE	
		PRINTED NAME SIGNATURE	PRINTED NAME SIGNATURE	000003

LABORATORY DELIVERABLES

		Check if Complete
1.	Cover page, Title page listing Lab Certification# facility name & address, & date of report	✓
2.	Table of Contents	✓
3.	Summary sheets listing analytical results for all targeted and non-targeted compounds	NA
4.	Summary Table cross-referencing field ID #'s vs. Lab ID #'s	✓
5.	Document bound, paginated and legible	✓
6.	Chain of Custody	✓
7.	Methodology Summary	✓
8.	Laboratory Chronicle and Holding Time check	✓
9.	Results submitted on a dry weight basis (if applicable)	NA
10.	Method Detection Limits	NA
11.	Lab certified by NJDEPE for parameters or appropriate category of parameters or a member of the USEPA CLP	✓
12.	Non-Conformance Summary	✓

Yvonne Berg — 2/21/95

Laboratory Manager or Environmental Consultant's Signature — Date

GC/MS ANALYSIS CONFORMANCE/NON-CONFORMANCE SUMMARY FORMAT

		No	Yes
1.	Chromatograms Labeled/Compounds Identified (Field Samples and Method Blanks)	___	✓
2.	GC/MS Tune Specifications		
	a. BFB Meet Criteria	___	✓
	b. DFTPP Meet Criteria	___	✓
3.	GC/MS Tune Frequency - Performed every 24 hours for 600 series and 12 hours for 8000 series.	___	✓
4.	GC/MS Calibration - Initial Calibration performed within 30 days before sample analysis and continuing calibration performed within 24 hours of sample analysis for 600 series and 12 hours for 8000 series	___	✓
5.	GC/MS Calibration Requirements		
	a. Calibration Check Compounds	___	✓
	b. System Performance Check Compounds	___	✓

6. Blank Contamination - If yes, list compounds and concentrations in each blank: Lm Jumalon 2/18/95

a. VOA fraction ~~VBlk 3 Methyl 2-pentanone, VBlk 4 Methyl 2-pentanone~~
b. ~~B/N Fraction~~ _____
c. Acid Fraction _____

~~VBlk 4 Methyl 2-pentanone 2.1 ppb Xylenes (total) 2 ppb~~

7. Surrogate Recoveries Meet Criteria ___ ✓ ~~✓~~ RH 1/11/95

If not met, list those compounds and their recoveries which fall outside the acceptable range:

a. VOA Fraction _____
b. B/N Fraction ~~See Form 2C~~ Lm 2/18/95
c. Acid Fraction _____

If not met, were the calculations checked and the results qualified as estimated? NA ✓ RH 1/11/95

8. Matrix Spike/Matrix Spike Duplicate Recoveries Meet Criteria (if not met, list these compounds and their recoveries which fall outside the acceptable range) ___ ✓

a. VOA Fraction Benzene (156)
b. B/N Fraction ~~See Reference Reports~~ Lm 2/18/95
c. Acid Fraction _____

9. Internal Standard Area/Retention Time Shift Meet Criteria ~~N/A~~

GC/MS ANALYSIS CONFORMANCE/NON-CONFORMANCE SUMMARY FORMAT (CONT.)

	No	Yes
10. Extraction Holding Time Met		NA

If not met, list number of days exceeded each sample: _____

11. Analysis Holding Time Met		✓

If not met, list number of days exceeded for each sample: _____

Additional Comments: _____

Laboratory Manager: [signature] Date 2/21/95

METAL ANALYSIS CONFORMANCE/NON-CONFORMANCE SUMMARY FORMAT

		No	Yes
1.	Calibration Summary Meet Criteria	___	✓
2.	ICP Interference Check Sample Results Summary Submitted (if applicable) / Meet Criteria	NA	___
3.	Serial Dilution Summary Submitted (if applicable) / Meet Criteria	NA	✓ (ES 2/9/95)
4.	Laboratory Control Sample Summary Submitted (if applicable) / Meet Criteria	___	✓
5.	Blank Contamination - If yes, list compounds concentrations in each blank:	✓	___
6.	Matrix Spike/Matrix Spike Duplicate Recoveries Meet Criteria (If not met, list those compounds and their recoveries which fall outside the acceptable range) SEE FORMS...	✓	___
7.	Extraction Holding Time Met	___	✓
	If not met, list number of days exceeded for each sample:		
8.	Analysis Holding Time Met	___	✓
	If not met, list number of days exceeded for each sample:		

Additional Comments: _____

Laboratory Manager: [signature] Date: 2/21/95

Laboratory Chronicle

Client Name: Aguilar Associates

Log In No.: 22633B

Date(s) of Sample Collection: 11/29/94

Date Received: 11/30/94

Sample ID: As per chain of custody

Organics Extraction:

1. Acids
2. Base/Neutrals
3. Pesticides/PCBs
4. Herbicides

Analysis:

12/1/94

1. Volatiles
2. Acids
3. Base/Neutrals
4. Pesticides/PCBs
5. Herbicides

Section Supervisor
Review & Approval

Digestion - 12/06/94 Analysis - 12/16/94, 12/19/94

Inorganics:

1. Metals
2. Cyanides
3. Phenols

Other Analysis:

Section Supervisor
Review & Approval

Quality Control Supervisor
Review & Approval

Dates are included for re-extractions and reanalysis.

NARRATIVE DISCUSSION
VOLATILES - 22633B
==========
Building 209

INTRODUCTION

This narrative covers the analysis of five (5) samples in accordance with the protocols based on NEI's SOP #703 and on EPA Method 624.

HOLDING TIMES

The analytical holding time for this analysis was met.

CALIBRATIONS

All required minimum RRFs and maximum % RSD initial calibration requirements have been met in accordance with the Method.

QC CHECK SAMPLE

All % recoveries in the QC check sample met the requirements as stipulated by the Method.

METHOD BLANK

The method blank associated with these samples met all method requirements.

SURROGATES

All surrogate recoveries met QC criteria.

MATRIX SPIKES

As requested, sample 296-6 was utilized for the MS/MSD. Spike recoveries fell with in QC limits, with the exception of 296-6MSD, which yielded a high Benzene recovery of 155.95 %. The form 3 is included with this report.

INTERNAL STANDARDS

Internal Standard area response/retention time summaries are not required.

SAMPLE COMMENTS

The TICs identified as "Unknown Siloxane" are most probably due to column degradation and not sample constituency.
NEI is reporting the results to our method detection limits (MDL's) rounded up to the nearest part per billion (ppb) in accordance with the guidance provided by NJDEP. These MDL's indicate that NEI did not detect andy compounds above these levels.
No further analytical problems were encountered.

NARRATIVE DISCUSSION
INORGANICS - 22603

All samples were analyzed as per the required protocol.

The matrix spike recovery was above control limits for sample 2044-3MS, thus resulting in the qualification of the sample results with "N". A second spike was performed on sample 296-6 and the recovery was within control limits.

The duplicate of sample 2044-3 showed no RPD, resulting in no additional sample qualification. A second duplicate was performed on sample 296-6. This second duplicate demonstrated a RPD of 8.2.

2/24/95

nytest environmental inc

I certify that this data package has been reviewed for the quality control and quality assurance measures for all analyzed methodologies.

Remo Gigante
Exec. Vice President

METHODOLOGY SUMMARY

AQUEOUS METHODOLOGIES:	REF 1	REF 2	REF 3	REF 5
BNA, Pesticides/PCB's Extraction		3510/3520		
AA/ICP Sample Preparation	200.7			
Furnace Sample Preparation	200.0			
Mercury Sample Preparation	245.1			
Hexavalent Chromium Sample Preparation	218.5			
Clean-Up		3610/3620/3630/ 3640/3660		
Organochlorine Pesticide and PCB's by Gas Chromatography			608	505
Herbicides by Gas Chromatography			362	515.1
Purgeable Organics by GC/MS			624	524.2
Base/Neutral, Acids by GC/MS			625	525
2,3,7,8-TCDD by GC/MS			613/625	
BTEX			602	502.2
EDB/DBCP by Microextraction				504.1

NON-AQUEOUS METHODOLOGIES:	
BNA, Pesticides/PCB's Extraction	3550
AA/ICP Sample Preparation	3050
Furnace Sample Preparation	3020/3030/3050
Mercury Sample Preparation	7471
Clean-Up	3610/3620/3630/ 3640/3660

GC, Gas Chromatography/Mass Spectrometry:	
Purgeable Organics	8240/8021
Base/Neutral and Acid Extractables	8270
Organophosphorus Pesticides	8140
Organochlorine Pesticide and PCB's by Gas Chromatography	8080
BTEX	8020
Halogenated Purgeable Organics	8010

METHODOLOGY SUMMARY

INDUCTIVELY COUPLED PLASMA (ICP):	REFERENCE 1	REFERENCE 2
Aluminum	200.7	6010
Antimony	200.7	6010
Barium	200.7	6010
Beryllium	200.7	6010
Cadmium	200.7	6010
Calcium	200.7	6010
Chromium	200.7	6010
Cobalt	200.7	6010
Copper	200.7	6010
Iron	200.7	6010
Lead	200.7	6010
Magnesium	200.7	6010
Manganese	200.7	6010
Molybdenum	200.7	6010
Nickel	200.7	6010
Potassium	200.7	6010
Silver	200.7	6010
Sodium	200.7	6010
Tin	200.7	6010
Titanium	200.7	6010
Vanadium	200.7	6010
Zinc	200.7	6010
FURNACE AA:		
Antimony	204.1	7041
Arsenic	206.2	7060
Lead	239.2	7421
Selenium	270.2	7740
Thallium	279.2	7841
Tin	282.2	
Vanadium	286.2	7911
Mercury	245.1	7470/7471
ICAP:		
Priority Pollutants	200.7	6010/7060/ 7470/7740
TAL Metals	200.7	6010/7060/ 7470/7740
RCRA Metals	200.7	6010/7060/ 7470/7740

METHODOLOGY SUMMARY

REFERENCES:

(1) USEPA-600/4-79-020, Methods for Chemical Analysis of Water and Waste

(2) USEPA SW 846, Test Methods for Evaluating Solid Waste, Third Edition

(3) Federal Register 40 CFR Part 136, Vol.49, No.209 Test Parameters for the Analysis of Pollutants

(4) Federal Register Vol.51, No.216 Friday, 11/7/86, pp.40643-40652

(5) Method for the Determination of Organic Compounds in Drinking Water, EPA 500/4-88/039, Dec. 1988

(6) Standard Method for Examination of Water and Wastewater, 15 Edition 1980

Method Qualifiers for Organic Non-CLP Methodologies

Q Qualifier - Specified entries and their meanings as follows:

U - Indicates compound was analyzed for but was not detected. The sample quantitation limit is corrected for dilutions and for the moisture content for soil samples. If a sample extract can not be concentrated to the protocol - specific volume, this fact is also accounted for in reporting the sample quantitation limit. The number is the minimum detected limits for the sample.

J - Indicates an estimated volume. The flag is used either when estimating concentration for tentatively identified compounds where a 1:1 response is assumed, or when the mass spectral data indicates the presence of a compound that meets the identification criteria but the result is less than the sample quantitation limit but greater than zero.

N - Indicates presumptive evidence of a compound. This flag is used only for tentatively identified compounds, where the identification is based on a mass spectral library search. It is applied to all TIC results. For generic characterization of a TIC, such as chlorinated hydrocarbon, the **N** code is not used.

B - This flag is used when the analyte is found in the analyte is found in the associated blank as well as the sample. It indicates possible/probable blank contamination and warns the data user to take appropriate action. This flag is used for a TIC as well as for a positively identified target compound.

E - This flag identifies compounds whose concentrations exceeded the calibration range of the GC/MS instrument for that specific analysis.

D - This flag identifies all compounds identified in an analysis at a secondary dilution factor.

A - This flag indicates that a TIC is a suspected aldol condensation product.

Method Qualifiers for Inorganics

FORM I-IN includes fields for three types of results qualifiers. These qualifiers must be completed as follows:

* C (Concentration) qualifier -- Enter "B" if the reported value was obtained from a reading that was less than the Contract Required Detection Limit (CRDL) but greater than or equal to the Instrument Detection Limit (IDL). If the analyte was analyzed for but not detected, a "U" must be entered.

* Q Qualifier -- Specified entries and their meanings are as follows :

E - The reported value is estimated because of the presence of interference.

M - Duplicate precision not met (CV > 20%).

N - Spiked sample recovery not within control limits.

S - The reported value was determined by Method of Standard Addition (MSA).

W - Post-digestion spike for Furnace AA analysis is out of control limits (85-115%), while sample absorbance is less than 50% of spike absorbance.

* - Duplicate analysis not within control limits.

\+ - Correlation Coefficient for MSA is less than 0.995.

Entering "S", "W" or "+" is mutually exclusive.

* M (Method) qualifier - enter:

- "P" for ICP
- "A" for Flame AA
- "F" for Furnace AA
- "CV" for Cold Vapor AA
- "AV" for Automated Cold Vapor AA
- "AS" for Semi-Automated Spectrophotometric
- "C" for Manual Spectrophotometric
- "T" for Titrimetric
- "NR" if the analyte is not required to be analyzed.

GC/MS Data

Volatile Data

1A

VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

296-FB

Lab Name: NYTEST ENV INC Contract: 9421415

Lab Code: NYTEST Case No.: 22633 SAS No.: SDG No.: ARMY3

Matrix: (soil/water) WATER Lab Sample ID: 2263307

Sample wt/vol: 5.0 (g/mL) ML Lab File ID: M1154.D

Level: (low/med) LOW Date Received: 11/30/94

% Moisture: not dec. _____ Date Analyzed: 12/01/94

Column: (pack/cap) CAP Dilution Factor: 1.0

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) UG/L	Q
74-87-3	Chloromethane	2	U
74-83-9	Bromomethane	1	U
75-01-4	Vinyl Chloride	1	U
75-00-3	Chloroethane	1	U
75-09-2	Methylene Chloride	3	U
75-35-4	1,1-Dichloroethene	2	U
75-34-3	1,1-Dichloroethane	1	U
67-66-3	Chloroform	1	U
107-06-2	1,2-Dichloroethane	1	U
71-55-6	1,1,1-Trichloroethane	1	U
56-23-5	Carbon Tetrachloride	2	U
75-27-4	Bromodichloromethane	1	U
78-87-5	1,2-Dichloropropane	1	U
10061-01-5	cis-1,3-Dichloropropene	1	U
79-01-6	Trichloroethene	2	U
124-48-1	Dibromochloromethane	1	U
79-00-5	1,1,2-Trichloroethane	1	U
71-43-2	Benzene	1	U
10061-02-6	trans-1,3-Dichloropropene	1	U
75-25-2	Bromoform	1	U
127-18-4	Tetrachloroethene	3	U
79-34-5	1,1,2,2-Tetrachloroethane	2	U
108-88-3	Toluene	2	U
108-90-7	Chlorobenzene	2	U
100-41-4	Ethylbenzene	2	U
1330-20-7	Xylene (total)	6	U
75-69-4	Trichloromonofluoromethane	2	U
107-02-8	Acrolein	20	U
107-13-1	Acrylonitrile	2	U
75-65-0	Tertiary Butyl Alcohol	100	U
1634-34-4	Methyl Tertiary Butyl Ether	1	U
541-73-1	1,3-Dichlorobenzene	2	U
106-46-7	1,4-Dichlorobenzene	2	U
95-50-1	1,2-Dichlorobenzene	2	U

1A

VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

296-FB

Lab Name: NYTEST ENV INC Contract: 9421415

Lab Code: NYTEST Case No.: 22633 SAS No.: SDG No.: ARMY3

Matrix: (soil/water) WATER Lab Sample ID: 2263307

Sample wt/vol: 5.0 (g/mL) ML Lab File ID: M1154.D

Level: (low/med) LOW Date Received: 11/30/94

% Moisture: not dec. _____ Date Analyzed: 12/01/94

Column: (pack/cap) CAP Dilution Factor: 1.0

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) UG/L	Q
110-75-8	2-Chloroethylvinyl Ether	4	U
156-60-5	Trans, 1,2-Dichloroethene	1	U

1E

EPA SAMPLE NO.

VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

296-FB

Lab Name: NYTEST ENV INC Contract: 9421415

Lab Code: NYTEST Case No.: 22633 SAS No.: SDG No.: ARMY3

Matrix: (soil/water) WATER Lab Sample ID: 2263307

Sample wt/vol: 5.0 (g/mL) ML Lab File ID: M1154.D

Level: (low/med) LOW Date Received: 11/30/94

% Moisture: not dec. _____ Date Analyzed: 12/01/94

Column: (pack/cap) CAP Dilution Factor: 1.0

Number TICs found: 2 CONCENTRATION UNITS: (ug/L or ug/Kg) UG/L

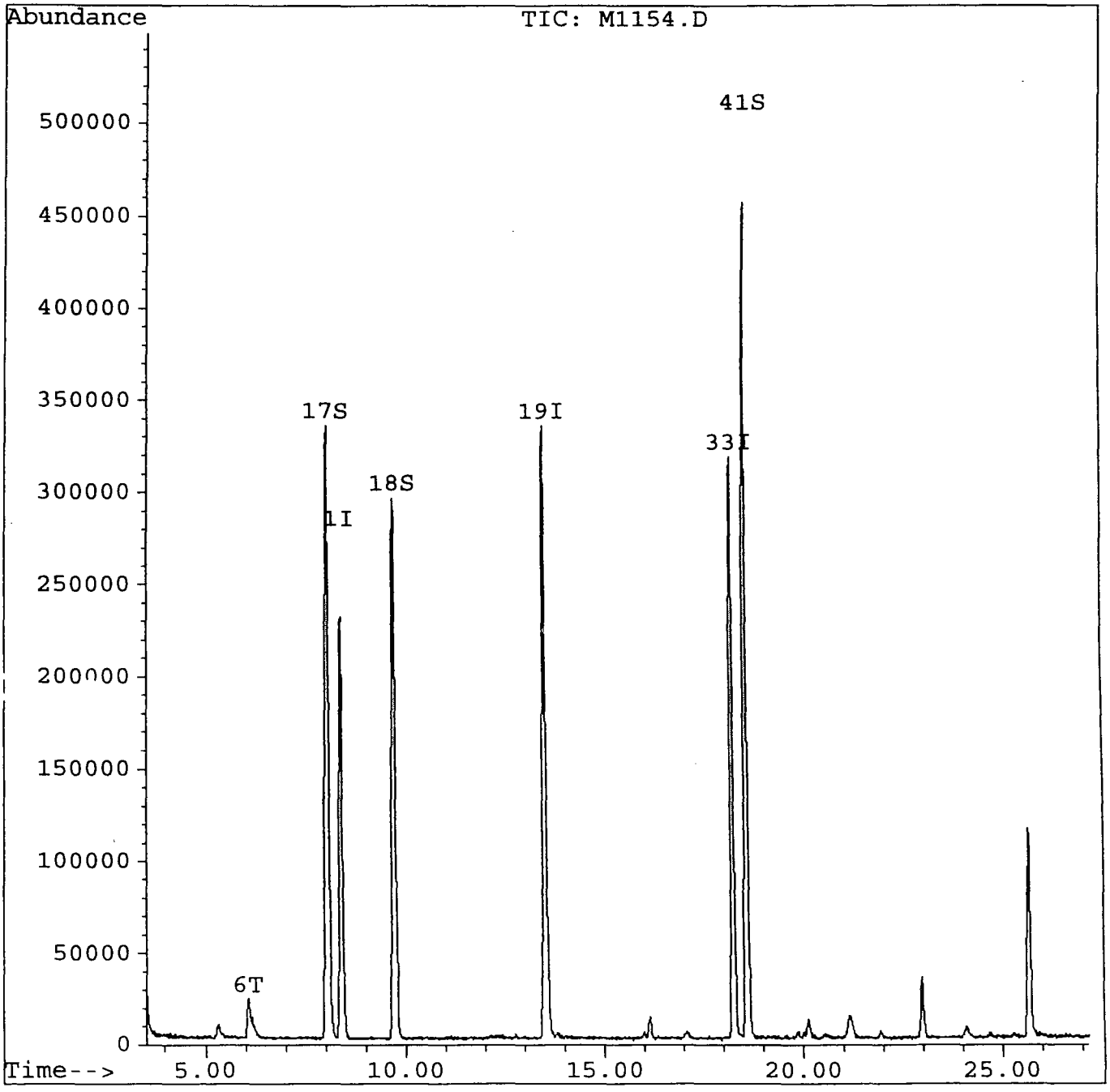
	CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1.		UNKNOWN HYDROCARBON	22.983	3	J
2.		UNKNOWN HYDROCARBON	25.671	12	J
3.					
4.					
5.					
6.					
7.					
8.					
9.					
10.					
11.					
12.					
13.					
14.					
15.					
16.					
17.					
18.					
19.					
20.					
21.					
22.					
23.					
24.					
25.					
26.					
27.					
28.					
29.					
30.					

Quantitation Report

Data File : C:\HPCHEM\1\DATA\120194\M1154.D
Acq Time : 1 Dec 94 13:48 pm
Sample : 2263307,296-FB,
Misc : 1,1,,,5,5,L,W,R11-30-94,
Quant Time: Dec 1 14:15 1994

Operator: VC
Inst : HPM
Multiplr: 1.00

Method : C:\HPCHEM\1\METHODS\P624.M
Title : VOA Standards for 5 point calibration
Last Update : Tue Nov 29 12:18:35 1994
Response via : Multiple Level Calibration



000022

Quantitation Report

Data File : C:\HPCHEM\1\DATA\120194\M1154.D
Acq Time : 1 Dec 94 13:48 pm
Sample : 2263307,296-FB,
Misc : 1,1,,,5,5,L,W,R11-30-94,
Quant Time: Dec 1 14:15 1994

Operator: VC
Inst : HPM
Multiplr: 1.00

Method : C:\HPCHEM\1\METHODS\P624.M
Title : VOA Standards for 5 point calibration
Last Update : Tue Nov 29 12:18:35 1994
Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) CI01 Bromochloromethane	8.39	128	141796	30.00	ug/l	-0.03
19) 2-Bromo-1-Chloropropane	13.52	77	505610	30.00	ug/l	-0.04
33) 1,4-Dichlorobutane	18.24	55	561804	30.00	ug/l	-0.04
System Monitoring Compounds						%Recovery
17) Pentafluorobenzene	8.05	168	867559	27.53	ug/l	91.78%
18) Fluorobenzene	9.72	96	748694	27.34	ug/l	91.14%
41) CS10 4-Bomofluorobenzene	18.59	95	435757	27.24	ug/l	90.80%
Target Compounds						Qvalue
6) C030 Methylene Chloride L	6 08	84	32041	2.46	ug/l	96

L COMPOUND BELOW MRL

WE 1/10/95

000023

(#) = qualifier out of range (m) = manual integration

1A

VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

296-TB

Lab Name: NYTEST ENV INC Contract: 9421415

Lab Code: NYTEST Case No.: 22633 SAS No.: SDG No.: ARMY3

Matrix: (soil/water) WATER Lab Sample ID: 2263308

Sample wt/vol: 5.0 (g/mL) ML Lab File ID: M1153.D

Level: (low/med) LOW Date Received: 11/30/94

% Moisture: not dec. _____ Date Analyzed: 12/01/94

Column: (pack/cap) CAP Dilution Factor: 1.0

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) UG/L	Q
74-87-3	Chloromethane	2	U
74-83-9	Bromomethane	1	U
75-01-4	Vinyl Chloride	1	U
75-00-3	Chloroethane	1	U
75-09-2	Methylene Chloride	3	
75-35-4	1,1-Dichloroethene	2	U
75-34-3	1,1-Dichloroethane	1	U
67-66-3	Chloroform	1	U
107-06-2	1,2-Dichloroethane	1	U
71-55-6	1,1,1-Trichloroethane	1	U
56-23-5	Carbon Tetrachloride	2	U
75-27-4	Bromodichloromethane	1	U
78-87-5	1,2-Dichloropropane	1	U
10061-01-5	cis-1,3-Dichloropropene	1	U
79-01-6	Trichloroethene	2	U
124-48-1	Dibromochloromethane	1	U
79-00-5	1,1,2-Trichloroethane	1	U
71-43-2	Benzene	1	U
10061-02-6	trans-1,3-Dichloropropene	1	U
75-25-2	Bromoform	1	U
127-18-4	Tetrachloroethene	3	U
79-34-5	1,1,2,2-Tetrachloroethane	2	U
108-88-3	Toluene	2	U
108-90-7	Chlorobenzene	2	U
100-41-4	Ethylbenzene	2	U
1330-20-7	Xylene (total)	6	U
75-69-4	Trichloromonofluoromethane	2	U
107-02-8	Acrolein	20	U
107-13-1	Acrylonitrile	2	U
75-65-0	Tertiary Butyl Alcohol	100	U
1634-34-4	Methyl Tertiary Butyl Ether	1	U
541-73-1	1,3-Dichlorobenzene	2	U
106-46-7	1,4-Dichlorobenzene	2	U
95-50-1	1,2-Dichlorobenzene	2	U

1A

VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

296-TB

Lab Name: NYTEST ENV INC Contract: 9421415

Lab Code: NYTEST Case No.: 22633 SAS No.: SDG No.: ARMY3

Matrix: (soil/water) WATER Lab Sample ID: 2263308

Sample wt/vol: 5.0 (g/mL) ML Lab File ID: M1153.D

Level: (low/med) LOW Date Received: 11/30/94

% Moisture: not dec. _____ Date Analyzed: 12/01/94

Column: (pack/cap) CAP Dilution Factor: 1.0

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) UG/L	Q
110-75-8	2-Chloroethylvinyl Ether	4	U
156-60-5	Trans, 1,2-Dichloroethene	1	U

000025

1E

EPA SAMPLE NO.

VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

296-TB

Lab Name: NYTEST ENV INC Contract: 9421415

Lab Code: NYTEST Case No.: 22633 SAS No.: SDG No.: ARMY3

Matrix: (soil/water) WATER Lab Sample ID: 2263308

Sample wt/vol: 5.0 (g/mL) ML Lab File ID: M1153.D

Level: (low/med) LOW Date Received: 11/30/94

% Moisture: not dec. _____ Date Analyzed: 12/01/94

Column: (pack/cap) CAP Dilution Factor: 1.0

Number TICs found: 1

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/L

	CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1.		UNKNOWN HYDROCARBON	25.661	13	J
2.					
3.					
4.					
5.					
6.					
7.					
8.					
9.					
10.					
11.					
12.					
13.					
14.					
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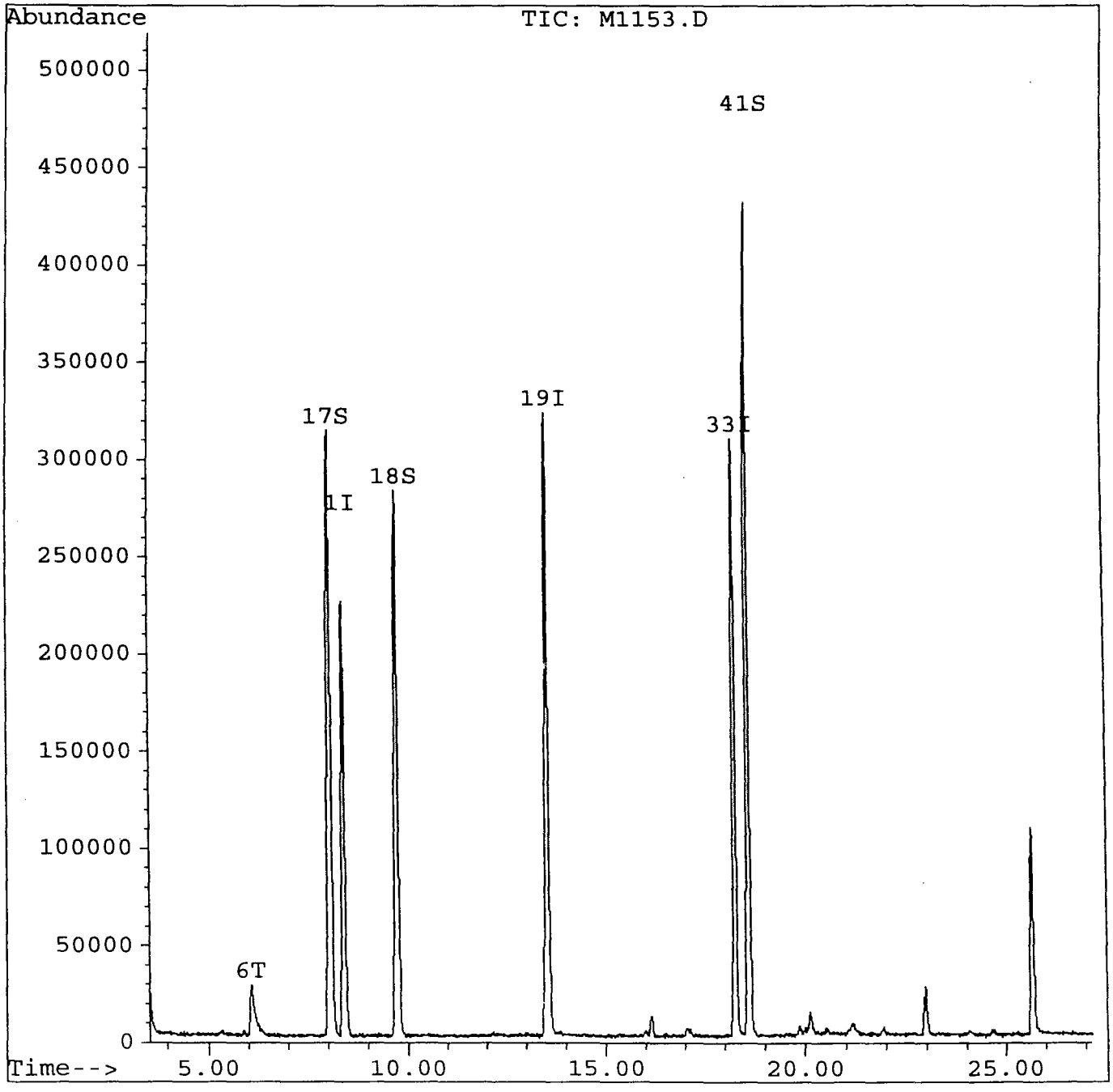
16.					
17.					
18.					
19.					
20.					
21.					
22.					
23.					
24.					
25.					
26.					
27.					
28.					
29.					
30.					

Quantitation Report

Data File : C:\HPCHEM\1\DATA\120194\M1153.D
Acq Time : 1 Dec 94 13:15 pm
Sample : 2263308,296-TB,
Misc : 1,1,,,5,5,L,W,R11-30-94,
Quant Time: Dec 1 13:43 1994

Operator: VC
Inst : HPM
Multiplr: 1.00

Method : C:\HPCHEM\1\METHODS\P624.M
Title : VOA Standards for 5 point calibration
Last Update : Tue Nov 29 12:18:35 1994
Response via : Multiple Level Calibration



000027

Quantitation Report

Data File : C:\HPCHEM\1\DATA\120194\M1153.D
Acq Time : 1 Dec 94 13:15 pm
Sample : 2263308,296-TB,
Misc : 1,1,,,5,5,L,W,R11-30-94,
Quant Time: Dec 1 13:43 1994

Operator: VC
Inst : HPM
Multiplr: 1.00

Method : C:\HPCHEM\1\METHODS\P624.M
Title : VOA Standards for 5 point calibration
Last Update : Tue Nov 29 12:18:35 1994
Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) CI01 Bromochloromethane	8.39	128	137091	30.00	ug/l	-0.03
19) 2-Bromo-1-Chloropropane	13.52	77	482268	30.00	ug/l	-0.04
33) 1,4-Dichlorobutane	18.24	55	539346	30.00	ug/l	-0.04
System Monitoring Compounds						%Recovery
17) Pentafluorobenzene	8.05	168	820233	26.93	ug/l	89.75%
18) Fluorobenzene	9.72	96	708576	26.76	ug/l	89.21%
41) CS10 4-Bomofluorobenzene	18.59	95	413779	26.94	ug/l	89.81%
Target Compounds						Qvalue
6) C030 Methylene Chloride	6.08	84	41773	3.32	ug/l #	87

000028

(#) = qualifier out of range (m) = manual integration

1A

VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

290-1

Lab Name: NYTEST ENV INC Contract: 9421415

Lab Code: NYTEST Case No.: 22633 SAS No.: SDG No.: ARMY3

Matrix: (soil/water) WATER Lab Sample ID: 2263312

Sample wt/vol: 5.0 (g/mL) ML Lab File ID: M1165.D

Level: (low/med) LOW Date Received: 11/30/94

% Moisture: not dec. _____ Date Analyzed: 12/01/94

Column: (pack/cap) CAP Dilution Factor: 1.0

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) UG/L	Q
74-87-3	Chloromethane	2	U
74-83-9	Bromomethane	1	U
75-01-4	Vinyl Chloride	1	U
75-00-3	Chloroethane	1	U
75-09-2	Methylene Chloride	3	U
75-35-4	1,1-Dichloroethene	2	U
75-34-3	1,1-Dichloroethane	1	U
67-66-3	Chloroform	1	U
107-06-2	1,2-Dichloroethane	1	U
71-55-6	1,1,1-Trichloroethane	1	U
56-23-5	Carbon Tetrachloride	2	U
75-27-4	Bromodichloromethane	1	U
78-87-5	1,2-Dichloropropane	1	U
10061-01-5	cis-1,3-Dichloropropene	1	U
79-01-6	Trichloroethene	2	U
124-48-1	Dibromochloromethane	1	U
79-00-5	1,1,2-Trichloroethane	1	U
71-43-2	Benzene	1	U
10061-02-6	trans-1,3-Dichloropropene	1	U
75-25-2	Bromoform	1	U
127-18-4	Tetrachloroethene	3	U
79-34-5	1,1,2,2-Tetrachloroethane	2	U
108-88-3	Toluene	2	U
108-90-7	Chlorobenzene	2	U
100-41-4	Ethylbenzene	2	U
1330-20-7	Xylene (total)	6	U
75-69-4	Trichloromonofluoromethane	2	U
107-02-8	Acrolein	20	U
107-13-1	Acrylonitrile	2	U
75-65-0	Tertiary Butyl Alcohol	100	U
1634-34-4	Methyl Tertiary Butyl Ether	1	U
541-73-1	1,3-Dichlorobenzene	2	U
106-46-7	1,4-Dichlorobenzene	2	U
95-50-1	1,2-Dichlorobenzene	2	U

1A

VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

290-1

Lab Name: NYTEST ENV INC Contract: 9421415

Lab Code: NYTEST Case No.: 22633 SAS No.: SDG No.: ARMY3

Matrix: (soil/water) WATER Lab Sample ID: 2263312

Sample wt/vol: 5.0 (g/mL) ML Lab File ID: M1165.D

Level: (low/med) LOW Date Received: 11/30/94

% Moisture: not dec. _____ Date Analyzed: 12/01/94

Column: (pack/cap) CAP Dilution Factor: 1.0

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) UG/L	Q
110-75-8	2-Chloroethylvinyl Ether	4	U
156-60-5	Trans, 1,2-Dichloroethene	1	U

1E

VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

290-1

Lab Name: NYTEST ENV INC Contract: 9421415

Lab Code: NYTEST Case No.: 22633 SAS No.: SDG No.: ARMY3

Matrix: (soil/water) WATER Lab Sample ID: 2263312

Sample wt/vol: 5.0 (g/mL) ML Lab File ID: M1165.D

Level: (low/med) LOW Date Received: 11/30/94

% Moisture: not dec. _____ Date Analyzed: 12/01/94

Column: (pack/cap) CAP Dilution Factor: 1.0

Number TICs found: 1

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/L

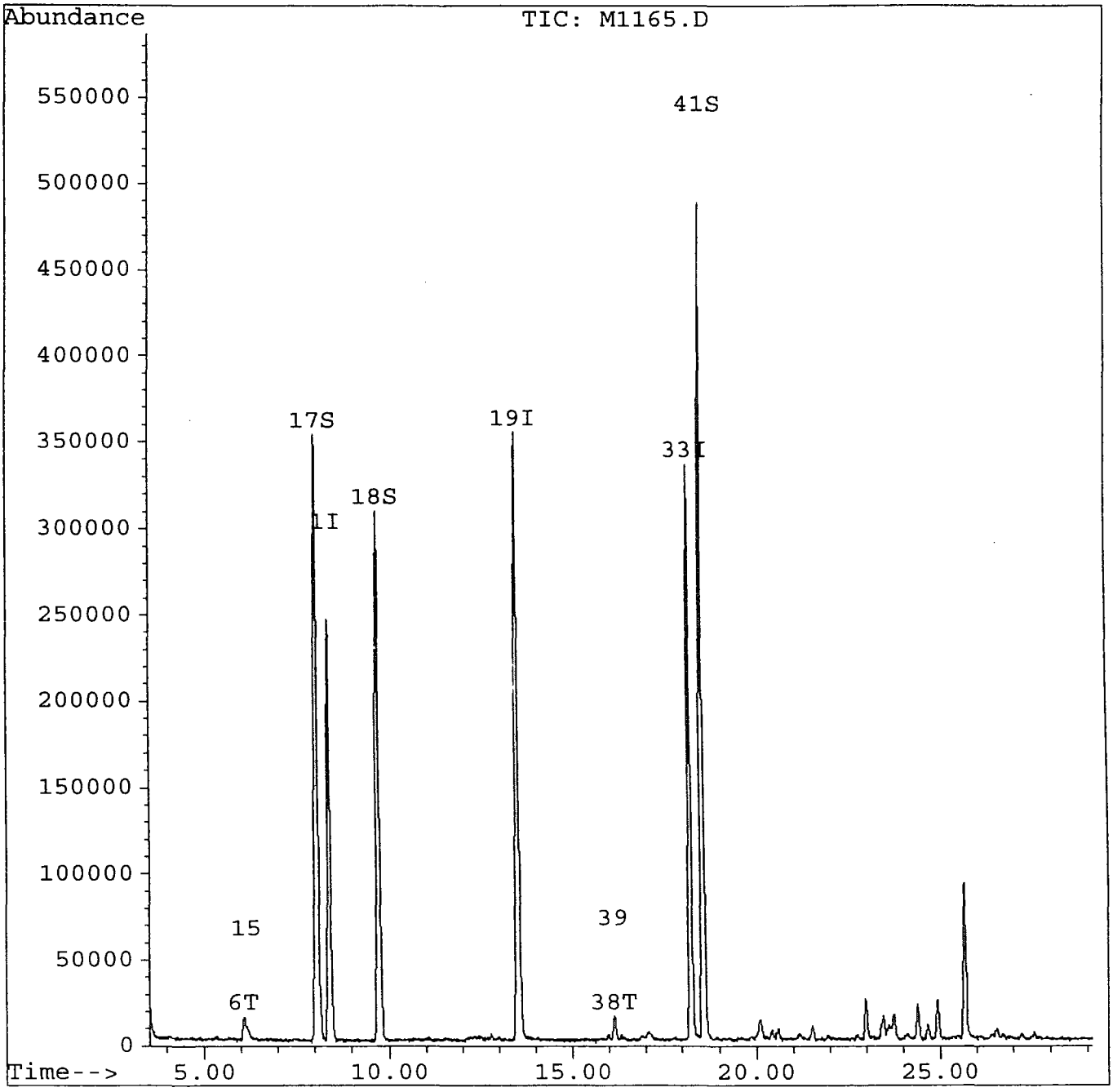
	CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1.		UNKNOWN	25.673	9	J
2.					
3.					
4.					
5.					
6.					
7.					
8.					
9.					
10.					
11.					
12.					
13.					
14.					
15.					
16.					
17.					
18.					
19.					
20.					
21.					
22.					
23.					
24.					
25.					
26.					
27.					
28.					
29.					
30.					

Quantitation Report

Data File : C:\HPCHEM\1\DATA\120194\M1165.D
Acq Time : 1 Dec 94 20:22 pm
Sample : 2263312,290-1,
Misc : 1,1,,,5,5,L,W,R11-30-94,
Quant Time: Dec 14 16:34 1994

Operator: VC
Inst : HPM
Multiplr: 1.00

Method : C:\HPCHEM\1\METHODS\P624.M
Title : VOA Standards for 5 point calibration
Last Update : Thu Dec 08 12:56:54 1994
Response via : Multiple Level Calibration



000032

Quantitation Report

Data File : C:\HPCHEM\1\DATA\120194\M1165.D
Acq Time : 1 Dec 94 20:22 pm
Sample : 2263312,290-1,
Misc : 1,1,,,5,5,L,W,R11-30-94,
Quant Time: Dec 14 16:34 1994

Operator: VC
Inst : HPM
Multiplr: 1.00

Method : C:\HPCHEM\1\METHODS\P624.M
Title : VOA Standards for 5 point calibration
Last Update : Thu Dec 08 12:56:54 1994
Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) CI01 Bromochloromethane	8.39	128	151134	30.00	ug/l	-0.03
19) 2-Bromo-1-Chloropropane	13.52	77	535265	30.00	ug/l	-0.04
33) 1,4-Dichlorobutane	18.24	55	593179	30.00	ug/l	-0.04
System Monitoring Compounds						%Recovery
17) Pentafluorobenzene	8.05	168	921870	27.45	ug/l	91.50%
18) Fluorobenzene	9.73	96	779231	26.70	ug/l	88.99%
41) CS10 4-Bomofluorobenzene	18.59	95	463642	27.45	ug/l	91.50%
Target Compounds						Qvalue
6) C030 Methylene Chloride L	6.08	84	13225	0.95	ug/l #	84
15) C176 Methyl Tertiary Butyl L	6.17	73	16451	0.94	ug/l	75
39) C250 M-P,Xylene L	16.14	106	13208	0.83	ug/l	90

L COMPOUND BELOW MDL WE 1/10/95

000033

(#) = qualifier out of range (m) = manual integration

1A

VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

VBLK39

Lab Name: NYTEST ENV INC Contract: 9421415

Lab Code: NYTEST Case No.: 22633 SAS No.: SDG No.: ARMY3

Matrix: (soil/water) WATER Lab Sample ID: VBLK39

Sample wt/vol: 5.0 (g/mL) ML Lab File ID: M1150.D

Level: (low/med) LOW Date Received: 00/00/00

% Moisture: not dec. _____ Date Analyzed: 12/01/94

Column: (pack/cap) CAP Dilution Factor: 1.0

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) UG/L	Q
74-87-3	Chloromethane	2	U
74-83-9	Bromomethane	1	U
75-01-4	Vinyl Chloride	1	U
75-00-3	Chloroethane	1	U
75-09-2	Methylene Chloride	3	U
75-35-4	1,1-Dichloroethene	2	U
75-34-3	1,1-Dichloroethane	1	U
67-66-3	Chloroform	1	U
107-06-2	1,2-Dichloroethane	1	U
71-55-6	1,1,1-Trichloroethane	1	U
56-23-5	Carbon Tetrachloride	2	U
75-27-4	Bromodichloromethane	1	U
78-87-5	1,2-Dichloropropane	1	U
10061-01-5	cis-1,3-Dichloropropene	1	U
79-01-6	Trichloroethene	2	U
124-48-1	Dibromochloromethane	1	U
79-00-5	1,1,2-Trichloroethane	1	U
71-43-2	Benzene	1	U
10061-02-6	trans-1,3-Dichloropropene	1	U
75-25-2	Bromoform	1	U
127-18-4	Tetrachloroethene	3	U
79-34-5	1,1,2,2-Tetrachloroethane	2	U
108-88-3	Toluene	2	U
108-90-7	Chlorobenzene	2	U
100-41-4	Ethylbenzene	2	U
1330-20-7	Xylene (total)	6	U
75-69-4	Trichloromonofluoromethane	2	U
107-02-8	Acrolein	20	U
107-13-1	Acrylonitrile	2	U
75-65-0	Tertiary Butyl Alcohol	100	U
1634-34-4	Methyl Tertiary Butyl Ether	1	U
541-73-1	1,3-Dichlorobenzene	2	U
106-46-7	1,4-Dichlorobenzene	2	U
95-50-1	1,2-Dichlorobenzene	2	U

1A

VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

VBLK39

Lab Name: NYTEST ENV INC Contract: 9421415

Lab Code: NYTEST Case No.: 22633 SAS No.: SDG No.: ARMY3

Matrix: (soil/water) WATER Lab Sample ID: VBLK39

Sample wt/vol: 5.0 (g/mL) ML Lab File ID: M1150.D

Level: (low/med) LOW Date Received: 00/00/00

% Moisture: not dec. _____ Date Analyzed: 12/01/94

Column: (pack/cap) CAP Dilution Factor: 1.0

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) UG/L	Q
110-75-8	2-Chloroethylvinyl Ether	4	U
156-60-5	Trans, 1,2-Dichloroethene	1	U

1E

VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

VBLK39

Lab Name: NYTEST ENV INC Contract: 9421415

Lab Code: NYTEST Case No.: 22633 SAS No.: SDG No.: ARMY3

Matrix: (soil/water) WATER Lab Sample ID: VBLK39

Sample wt/vol: 5.0 (g/mL) ML Lab File ID: M1150.D

Level: (low/med) LOW Date Received: 00/00/00

% Moisture: not dec. _____ Date Analyzed: 12/01/94

Column: (pack/cap) CAP Dilution Factor: 1.0

Number TICs found: 0 CONCENTRATION UNITS: (ug/L or ug/Kg) UG/L

	CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1.					
2.					
3.					
4.					
5.					
6.					
7.					
8.					
9.					
10.					
11.					
12.					
13.					
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23.					
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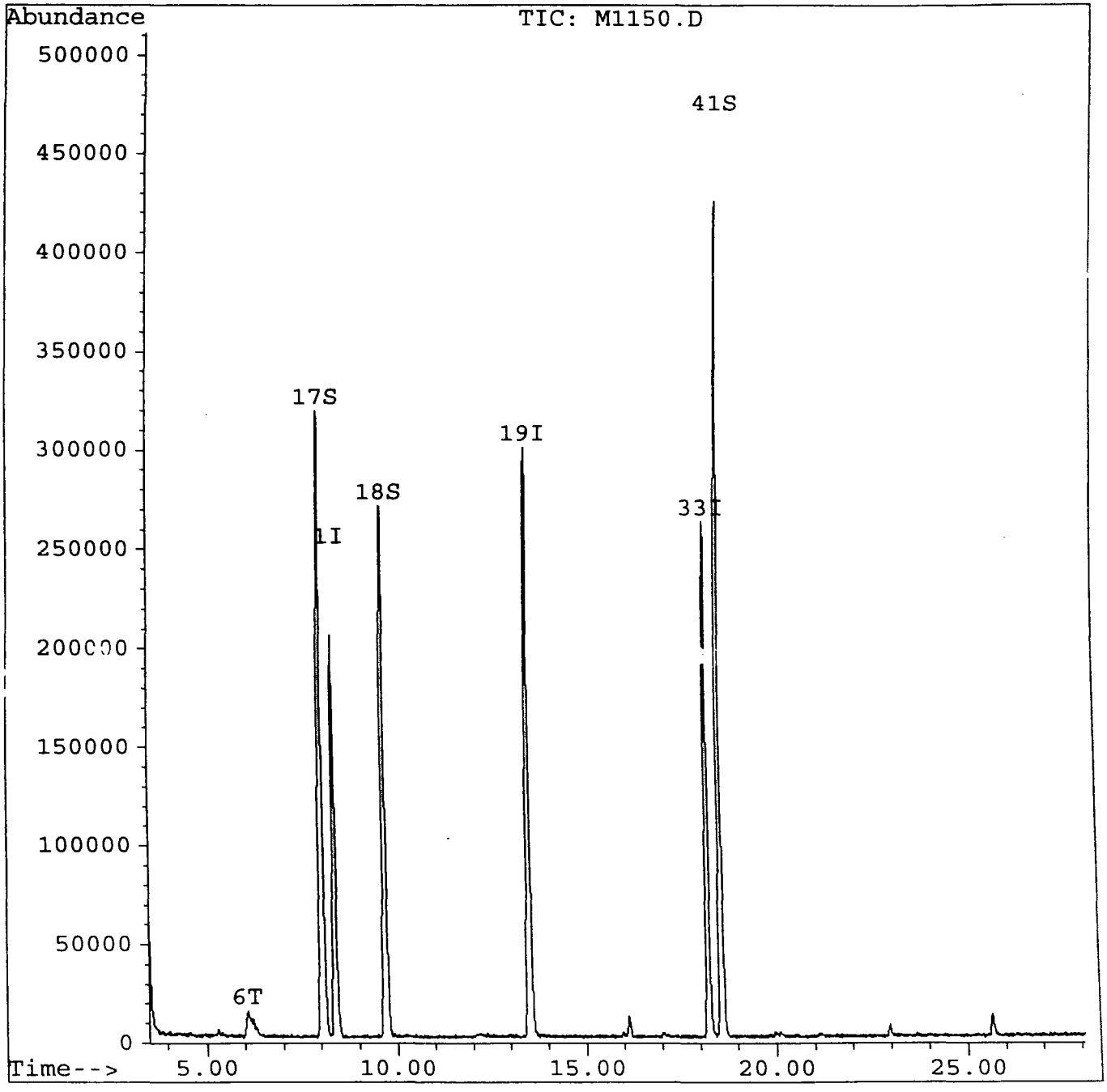
26.					
27.					
28.					
29.					
30.					

Quantitation Report

Data File : C:\HPCHEM\1\DATA\120194\M1150.D
Acq Time : 1 Dec 94 11:30 am
Sample : VBLK39,VBLK39,
Misc : 1,,,,,5,5,L,W,
Quant Time: Dec 1 12:00 1994

Operator: VC
Inst : HPM
Multiplr: 1.00

Method : C:\HPCHEM\1\METHODS\P624.M
Title : VOA Standards for 5 point calibration
Last Update : Tue Nov 29 12:18:35 1994
Response via : Multiple Level Calibration



000037

Quantitation Report

Data File : C:\HPCHEM\1\DATA\120194\M1150.D
Acq Time : 1 Dec 94 11:30 am
Sample : VBLK39,VBLK39,
Misc : 1,,,,,5,5,L,W,
Quant Time: Dec 1 12:00 1994

Operator: VC
Inst : HPM
Multiplr: 1.00

Method : C:\HPCHEM\1\METHODS\P624.M
Title : VOA Standards for 5 point calibration
Last Update : Tue Nov 29 12:18:35 1994
Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) CI01 Bromochloromethane	8.38	128	126949	30.00	ug/l	-0.05
19) 2-Bromo-1-Chloropropane	13.50	77	462593	30.00	ug/l	-0.06
33) 1,4-Dichlorobutane	18.23	55	462734	30.00	ug/l	-0.05
System Monitoring Compounds						%Recovery
17) Pentafluorobenzene	8.03	168	836157	29.64	ug/l	98.81%
18) Fluorobenzene	9.70	96	694203	28.32	ug/l	94.39%
41) CS10 4-Bomofluorobenzene	18.57	95	404891	30.73	ug/l	102.43%
Target Compounds						Qvalue
6) C030 Methylene Chloride 2	6.07	84	28277	2.42	ug/l	95

2 COMPOUNDS BELOW MDL

WE 1/10/95

000038

(#) = qualifier out of range (m) = manual integration

WATER VOLATILE SURROGATE COMPOUND RECOVERY

Lab Name: NYTEST ENV INC Contract: 9421415

Lab Code: NYTEST Case No.: 22633 SAS No.: SDG No.: ARMY3

	EPA SAMPLE NO.	SUR1 (FB) #	SUR2 (BFB)#	SUR3 (PFB)#	OTHER	TOT OUT
01	QCCHECK	88	93	89		0
02	VBLK39	94	102	99		0
03	2044-TB	92	99	94		0
04	296-TB	89	90	90		0
05	296-FB	91	91	92		0
06	2044-FB	94	94	94		0
07	296-6	92	93	92		0
08	296-6MS	91	89	92		0
09	296-1	91	93	91		0
10	296-3	90	93	92		0
11	296-2	89	88	91		0
12	296-7	89	91	90		0
13	296-8	92	93	94		0
14	296-DUP	90	94	92		0
15	290-1	89	92	92		0
16	108-1	88	91	90		0
17	108-2	91	94	93		0
18	108-3	89	92	92		0
19	750-1	89	91	91		0
20	750-3	88	91	89		0
21	750-4	90	92	92		0
22	2562-2	91	96	94		0
23	2562-3	90	94	92		0
24	2044-1	86	92	91		0
25	2044-2	86	81	86		0
26	2044-3	93	109	97		0
27	2044-DUP	92	101	95		0
28	2562-1	93	96	95		0
29	QCCHECK1	88	89	90		0
30	VBLK41	95	93	97		0

	QC LIMITS
SUR1 (FB) = Fluorobenzene	(70-140)
SUR2 (BFB) = Bromofluorobenzene	(60-150)
SUR3 (PFB) = Pentafluorobenzene	(70-140)

Column to be used to flag recovery values

* Values outside of contract required QC limits

D Surrogates diluted out

nytest

RECOVERY REPORT

Client Name:
Sample Matrix: LIQUID
Lab Smp Id: 2263310
Level: LOW
Data Type: MS DATA
SpikeList File: QCMS.spk
Method File: /chem/HPM.i/22633.b/624.m
Misc Info:

Client SDG: ARMY3
Fraction: VOA
Client Smp ID: 296-6MS
Operator: VC
SampleType: MS
Quant Type: ISTD

SPIKE	COMPOUND	CONC ADDED ug/l	CONC RECOVERED ug/l	% RECOVERED	LIMITS
1	Chloromethane	20	20	98.48	0-2[illegible]
4	Bromomethane	20	20	98.48	0-242
5	Vinyl Chloride	20	16	81.56	0-251
6	Chloroethane	20	19	93.87	14-230
7	Methylene Chloride	20	13	66.06	0-221
10	1,1-Dichloroethene	20	19	95.79	0-234
12	1,1-Dichloroethane	20	19	97.02	59-155
14	Trans, 1,2-Dichlor	20	19	97.24	54-156
15	Chloroform	20	20	98.97	51-138
16	1,2-Dichloroethane	20	19	95.69	49-155
19	Trichloromonofluor	20	20	101.35	17-181
24	1,1,1-Trichloroeth	20	19	93.48	52-162
25	Carbon Tetrachlori	20	18	91.23	70-140
27	Bromodichlorometha	20	18	91.55	35-155
28	1,2-Dichloropropan	20	18	90.83	0-210
29	cis-1,3-Dichloropr	20	18	92.90	0-227
30	Trichloroethene	20	18	88.48	71-157
31	Dibromochlorometha	20	18	89.02	53-149
32	1,1,2-Trichloroeth	20	18	90.41	52-150
34	Benzene	20	28	142.29	37-151
35	trans-1,3-Dichloro	20	18	92.74	17-183
37	Bromoform	20	18	91.99	45-169
38	2-Chloroethylvinyl	20	18	90.80	0-305
41	Tetrachloroethene	20	18	92.25	64-148
43	1,1,2,2-Tetrachlor	20	18	92.11	46-157
45	Toluene	20	19	95.44	47-150
46	Chlorobenzene	20	18	90.11	37-160
47	Ethylbenzene	20	26	130.08	37-162
53	1,3-Dichlorobenzen	20	20	99.34	59-156
54	1,4-Dichlorobenzen	20	21	104.51	18-190
55	1,2-Dichlorobenzen	20	21	103.70	18-190

nytest

RECOVERY REPORT

Client Name:
Sample Matrix: LIQUID
Lab Smp Id: 2263310
Level: LOW
Data Type: MS DATA
SpikeList File: QCMS.spk
Method File: /chem/HPM.i/22633.b/624.m
Misc Info:

Client SDG: ARMY3
Fraction: VOA
Client Smp ID: 296-6MS
Operator: VC
SampleType: MS
Quant Type: ISTD

SURROGATE COMPOUND	CONC ADDED ug/l	CONC RECOVERED ug/l	% RECOVERED	LIMITS
$ 33 Pentafluorobenzene	30	28	91.98	70-140
$ 44 Fluorobenzene	30	27	91.41	70-140
$ 48 Bromofluorobenzene	30	26	88.64	60-150

nytest

RECOVERY REPORT

Client Name:
Sample Matrix: LIQUID
Lab Smp Id: 2263311
Level: LOW
Data Type: MS DATA
SpikeList File: QCMS.spk
Method File: /chem/HPM.i/22633.b/624.m
Misc Info:

Client SDG: ARMY3
Fraction: VOA
Client Smp ID: ~~MW~~ 2W6-6MSD
Operator: VC
SampleType: MSD
Quant Type: ISTD

emt Summary 2/18/95

SURROGATE COMPOUND	CONC ADDED ug/l	CONC RECOVERED ug/l	% RECOVERED	LIMITS
$ 33 Pentafluorobenzene	30	27	90.25	70-140
$ 44 Fluorobenzene	30	27	89.48	70-140
$ 48 Bromofluorobenzene	30	27	88.92	60-150

4A

VOLATILE METHOD BLANK SUMMARY

EPA SAMPLE NO.

VBLK39

Lab Name: NYTEST ENV INC Contract: 9421415

Lab Code: NYTEST Case No.: 22633 SAS No.: SDG No.: ARMY3

Lab File ID: M1150.D Lab Sample ID: VBLK39

Date Analyzed: 12/01/94 Time Analyzed: 1130

Matrix: (soil/water) WATER Level: (low/med) LOW

Instrument ID: HPM

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES, MS AND MSD

	EPA SAMPLE NO.	LAB SAMPLE ID	LAB FILE ID	TIME ANALYZED
01	QCCHECK	QCCHECK	M1149.D	1040
02	2044-TB	2263327	M1151.D	1210
03	296-TB	2263308	M1153.D	1315
04	296-FB	2263307	M1154.D	1348
05	2044-FB	2263325	M1155.D	1430
06	296-6	2263304	M1156.D	1503
07	296-6MS	2263310	M1157.D	1536
08	296-1	2263301	M1159.D	1655
09	296-3	2263303	M1160.D	1730
10	296-2	2263302	M1161.D	1804
11	296-7	2263305	M1162.D	1839
12	296-8	2263306	M1163.D	1913
13	296-DUP	2263309	M1164.D	1948
14	290-1	2263312	M1165.D	2022
15	108-1	2263313	M1166.D	2057
16	108-2	2263314	M1167.D	2131
17	108-3	2263315	M1168.D	2205
18	750-1	2263316	M1169.D	2240
19	750-3	2263317	M1170.D	2314
20	750-4	2263318	M1171.D	2348
21	2562-2	2263320	M1173.D	0057
22	2562-3	2263321	M1174.D	0131
23	2044-1	2263322	M1175.D	0206
24	2044-2	2263323	M1176.D	0240
25	2044-3	2263324	M1177.D	0856
26	2044-DUP	2263326	M1178.D	0930
27	2562-1	2263319	M1179.D	1010
28				
29				
30				

COMMENTS:

5A

VOLATILE ORGANIC INSTRUMENT PERFORMANCE CHECK
BROMOFLUOROBENZENE (BFB)

Lab Name: NYTEST ENV INC Contract: 9421415

Lab Code: NYTEST Case No.: 22633 SAS No.: SDG No.: ARMY3

Lab File ID: M1098.D BFB Injection Date: 11/28/94

Instrument ID: HPM BFB Injection Time: 1706

Matrix: (soil/water) WATER Level: (low/med) LOW Column: (pack/cap) CAP

m/e	ION ABUNDANCE CRITERIA	% RELATIVE ABUNDANCE
50	15.0 - 40.0% of mass 95	24.7
75	30.0 - 60.0% of mass 95	46.6
95	Base peak, 100% relative abundance	100.0
96	5.0 - 9.0% of mass 95	6.3
173	Less than 2.0% of mass 174	0.0 (0.0)1
174	Greater than 50.0% of mass 95	95.9
175	5.0 - 9.0% of mass 174	7.3 (7.6)1
176	Greater than 95.0%, but less than 101.0% of mass 174	93.3 (97.2)1
177	5.0 - 9.0% of mass 176	6.1 (6.5)2

1-Value is % mass 174 2-Value is % mass 176

THIS CHECK APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS, AND STANDARDS:

	EPA SAMPLE NO.	LAB SAMPLE ID	LAB FILE ID	DATE ANALYZED	TIME ANALYZED
01	VSTD005	VSTD005	M1100.D	11/28/94	1755
02	VSTD020	VSTD020	M1102.D	11/28/94	1912
03	VSTD050	VSTD050	M1103.D	11/28/94	1950
04	VSTD100	VSTD100	M1105.D	11/28/94	2108
05	VSTD200	VSTD200	M1107.D	11/28/94	2225
06					
07					
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5A

VOLATILE ORGANIC INSTRUMENT PERFORMANCE CHECK
BROMOFLUOROBENZENE (BFB)

Lab Name: NYTEST ENV INC Contract: 9421415

Lab Code: NYTEST Case No.: 22633 SAS No.: SDG No.: ARMY3

Lab File ID: M1148.D BFB Injection Date: 12/01/94

Instrument ID: HPM BFB Injection Time: 1029

Matrix:(soil/water) WATER Level:(low/med) LOW Column:(pack/cap) CAP

m/e	ION ABUNDANCE CRITERIA	% RELATIVE ABUNDANCE
50	15.0 - 40.0% of mass 95	25.6
75	30.0 - 60.0% of mass 95	44.7
95	Base peak, 100% relative abundance	100.0
96	5.0 - 9.0% of mass 95	6.6
173	Less than 2.0% of mass 174	0.0 (0.0)1
174	Greater than 50.0% of mass 95	92.3
175	5.0 - 9.0% of mass 174	7.4 (8.0)1
176	Greater than 95.0%, but less than 101.0% of mass 174	90.4 (98.0)1
177	5.0 - 9.0% of mass 176	5.8 (6.4)2

1-Value is % mass 174 2-Value is % mass 176

THIS CHECK APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS, AND STANDARDS:

	EPA SAMPLE NO.	LAB SAMPLE ID	LAB FILE ID	DATE ANALYZED	TIME ANALYZED
01	QCCHECK	QCCHECK	M1149.D	12/01/94	1040
02	VBLK39	VBLK39	M1150.D	12/01/94	1130
03	2044-TB	2263327	M1151.D	12/01/94	1210
04	296-TB	2263308	M1153.D	12/01/94	1315
05	296-FB	2263307	M1154.D	12/01/94	1348
06	2044-FB	2263325	M1155.D	12/01/94	1430
07	296-6	2263304	M1156.D	12/01/94	1503
08	296-6MS	2263310	M1157.D	12/01/94	1536
09	296-1	2263301	M1159.D	12/01/94	1655
10	296-3	2263303	M1160.D	12/01/94	1730
11	296-2	2263302	M1161.D	12/01/94	1804
12	296-7	2263305	M1162.D	12/01/94	1839
13	296-8	2263306	M1163.D	12/01/94	1913
14	296-DUP	2263309	M1164.D	12/01/94	1948
15	290-1	2263312	M1165.D	12/01/94	2022
16	108-1	2263313	M1166.D	12/01/94	2057
17	108-2	2263314	M1167.D	12/01/94	2131
18	108-3	2263315	M1168.D	12/01/94	2205
19	750-1	2263316	M1169.D	12/01/94	2240
20	750-3	2263317	M1170.D	12/01/94	2314
21	750-4	2263318	M1171.D	12/01/94	2348
22	2562-2	2263320	M1173.D	12/02/94	0057

5A

VOLATILE ORGANIC INSTRUMENT PERFORMANCE CHECK
BROMOFLUOROBENZENE (BFB)

Lab Name: NYTEST ENV INC Contract: 9421415

Lab Code: NYTEST Case No.: 22633 SAS No.: SDG No.: ARMY3

Lab File ID: M1148.D BFB Injection Date: 12/01/94

Instrument ID: HPM BFB Injection Time: 1029

Matrix:(soil/water) WATER Level:(low/med) LOW Column:(pack/cap) CAP

m/e	ION ABUNDANCE CRITERIA	% RELATIVE ABUNDANCE
50	15.0 - 40.0% of mass 95	25.6
75	30.0 - 60.0% of mass 95	44.7
95	Base peak, 100% relative abundance	100.0
96	5.0 - 9.0% of mass 95	6.6
173	Less than 2.0% of mass 174	0.0 (0.0)1
174	Greater than 50.0% of mass 95	92.3
175	5.0 - 9.0% of mass 174	7.4 (8.0)1
176	Greater than 95.0%, but less than 101.0% of mass 174	90.4 (98.0)1
177	5.0 - 9.0% of mass 176	5.8 (6.4)2

1-Value is % mass 174 2-Value is % mass 176

THIS CHECK APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS, AND STANDARDS:

	EPA SAMPLE NO.	LAB SAMPLE ID	LAB FILE ID	DATE ANALYZED	TIME ANALYZED
01	2562-3	2263321	M1174.D	12/02/94	0131
02	2044-1	2263322	M1175.D	12/02/94	0206
03	2044-2	2263323	M1176.D	12/02/94	0240
04	2044-3	2263324	M1177.D	12/02/94	0856
05	2044-DUP	2263326	M1178.D	12/02/94	0930
06	2562-1	2263319	M1179 D	12/02/94	1010
07					
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6A

VOLATILE ORGANICS INITIAL CALIBRATION DATA

Lab Name: NYTEST ENV INC Contract: 9421415

Lab Code: NYTEST Case No.: 22633 SAS No.: SDG No.: ARMY3

Instrument ID: HPM Calibration Date(s): 11/28/94

Matrix:(soil/water) WATER Level:(low/med) LOW Column:(pack/cap) CAP

Max $\overline{\%RSD}$ for CCC(*) = 35.0%

LAB FILE ID: RRF005 =M1100.D RRF010 =M1102.D
RRF030=M1103.D RRF050=M1105.D RRF200=M1107.D

COMPOUND	RRF005	RRF010	RRF030	RRF050	RRF200	$\overline{RRF}$	% RSD
Chloromethane	1.236	1.732	1.432	1.462	1.359	1.444	12.7
Bromomethane	1.418	1.713	1.276	1.472	1.357	1.447	11.4
Vinyl Chloride	1.526	2.017	1.598	1.824	1.705	1.734	11.2
Chloroethane	0.911	1.227	0.942	1.102	1.040	1.044	12.2
Methylene Chloride	3.837	3.259	1.863	2.065		2.756	34.4
1,1-Dichloroethene	1.465	1.880	1.456	1.712	1.565	1.616	11.2
1,1-Dichloroethane	3.420	4.234	3.262	3.820	3.362	3.620	11.2
Chloroform	3.209	3.968	3.002	3.518	3.267	3.393	10.9
1,2-Dichloroethane	2.329	2.835	2.179	2.590	2.346	2.456	10.5
1,1,1-Trichloroethane	0.757	0.927	0.724	0.861	0.750	0.804	10.7
Carbon Tetrachloride	0.721	0.853	0.699	0.831	0.725	0.766	9.2
Bromodichloromethane	0.862	1.054	0.834	0.988	0.879	0.923	10.1
1,2-Dichloropropane	0.683	0.817	0.631	0.752	0.665	0.710	10.5
cis-1,3-Dichloropropene	0.858	1.041	0.820	0.960	0.836	0.903	10.5
Trichloroethene	0.638	0.770	0.606	0.701	0.633	0.670	9.9
Dibromochloromethane	0.774	0.985	0.783	0.934	0.826	0.860	10.9
1,1,2-Trichloroethane	0.540	0.664	0.516	0.606	0.531	0.571	10.9
Benzene	1.561	1.818	1.424	1.679	1.461	1.589	10.2
trans-1,3-Dichloropropene	0.745	0.901	0.711	0.835	0.731	0.784	10.2
Bromoform	0.715	0.942	0.779	0.920	0.812	0.834	11.5
Tetrachloroethene	0.698	0.780	0.604	0.708	0.632	0.684	10.1
1,1,2,2-Tetrachloroethane	0.827	0.999	0.768	0.906	0.761	0.851	11.9
Toluene	1.760	2.045	1.583	1.866	1.632	1.777	10.5
Chlorobenzene	1.209	1.365	1.042	1.226	1.093	1.187	10.6
Ethylbenzene	0.625	0.708	0.526	0.625	0.571	0.611	11.3
Xylene (total)	0.720	0.819	0.622	0.743	0.671	0.715	10.4
Trichloromonofluoromethane	2.566	3.298	2.535	2.927	2.714	2.808	11.2
Acrolein		0.109	0.121	0.047	0.106	0.096	34.9
Acrylonitrile	1.220	1.556	1.306	1.444	1.237	1.353	10.6
Tertiary Butyl Alcohol	0.000	0.204	0.165	0.165		0.178	12.6
Methyl Tertiary Butyl Ether	3.406	4.306	3.105	3.687	2.875	3.476	16.0
1,3-Dichlorobenzene	1.260	1.448	1.108	1.314	1.178	1.261	10.4
1,4-Dichlorobenzene	1.296	1.504	1.147	1.371	1.227	1.309	10.4
1,2-Dichlorobenzene	1.150	1.332	1.013	1.213	1.084	1.159	10.6
2-Chloroethylvinyl Ether	0.683	0.817	0.631	0.752	0.666	0.710	10.5
Trans, 1,2-Dichloroethene	1.574	1.988	1.517	1.798	1.596	1.695	11.5
Fluorobenzene	5.498	6.761	5.100	5.999	5.611	5.794	10.8

6A

VOLATILE ORGANICS INITIAL CALIBRATION DATA

Lab Name: NYTEST ENV INC Contract: 9421415

Lab Code: NYTEST Case No.: 22633 SAS No.: SDG No.: ARMY3

Instrument ID: HPM Calibration Date(s): 11/28/94

Matrix:(soil/water) WATER Level:(low/med) LOW Column:(pack/cap) CAP

Max $\overline{\%RSD}$ for CCC(*) = 35.0%

LAB FILE ID: RRF005 =M1100.D RRF010 =M1102.D
RRF030=M1103.D RRF050=M1105.D RRF200=M1107.D

COMPOUND	RRF005	RRF010	RRF030	RRF050	RRF200	$\overline{RRF}$	% RSD
Bromofluorobenzene	0.843	0.979	0.744	0.890	0.816	0.854	10.2
Pentafluorobenzene	6.449	7.843	5.883	6.876	6.279	6.666	11.2

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RECOVERY REPORT

Client Name:
Sample Matrix: LIQUID
Lab Smp Id: QCCHECK
Level: LOW
Data Type: MS DATA
SpikeList File: QCCHK.spk
Method File: /chem/HPM.i/22633.b/624.m
Misc Info:

Client SDG: ARMY3
Fraction: VOA
Client Smp ID: QCCHECK
Operator: VC
SampleType: MS
Quant Type: ISTD

SPIKE	COMPOUND	CONC ADDED ug/l	CONC RECOVERED ug/l	% RECOVERED	LIMITS
1	Chloromethane	20	20	99.67	0-204
4	Bromomethane	20	20	100.70	14-186
5	Vinyl Chloride	20	16	81.08	4-196
6	Chloroethane	20	19	94.22	38-162
7	Methylene Chloride	20	15	77.31	60-140
10	1,1-Dichloroethene	20	19	95.82	50-150
12	1,1-Dichloroethane	20	19	95.98	72-128
14	Trans, 1,2-Dichlor	20	20	97.58	70-130
15	Chloroform	20	20	98.98	68-132
16	1,2-Dichloroethane	20	19	94.03	68-132
19	Trichloromonofluor	20	20	101.55	48-152
24	1,1,1-Trichloroeth	20	20	99.39	75-125
25	Carbon Tetrachlori	20	20	98.47	73-127
27	Bromodichlorometha	20	19	95.92	66-134
28	1,2-Dichloropropan	20	19	95.91	34-166
29	cis-1,3-Dichloropr	20	20	97.81	24-176
30	Trichloroethene	20	19	95.80	66-134
31	Dibromochlorometha	20	19	95.12	68-132
32	1,1,2-Trichloroeth	20	19	94.43	71-129
34	Benzene	20	19	96.01	64-136
35	trans-1,3-Dichloro	20	19	96.86	50-150
37	Bromoform	20	19	96.94	71-129
38	2-Chloroethylvinyl	20	19	95.87	0-224
41	Tetrachloroethene	20	21	103.50	74-126
43	1,1,2,2-Tetrachlor	20	20	101.93	60-140
45	Toluene	20	20	98.58	74-126
46	Chlorobenzene	20	20	102.22	66-134
47	Ethylbenzene	20	20	100.60	59-141
53	1,3-Dichlorobenzen	20	22	109.06	73-127
54	1,4-Dichlorobenzen	20	23	116.98	63-137
55	1,2-Dichlorobenzen	20	23	115.29	63-127

nytest

RECOVERY REPORT

Client Name:
Sample Matrix: LIQUID
Lab Smp Id: QCCHECK
Level: LOW
Data Type: MS DATA
SpikeList File: QCCHK.spk
Method File: /chem/HPM.i/22633.b/624.m
Misc Info:

Client SDG: ARMY3
Fraction: VOA
Client Smp ID: QCCHECK
Operator: VC
SampleType: MS
Quant Type: ISTD

SURROGATE COMPOUND	CONC ADDED ug/l	CONC RECOVERED ug/l	% RECOVERED	LIMITS
$ 33 Pentafluorobenzene	30	27	89.43	70-140
$ 44 Fluorobenzene	30	26	87.62	70-140
$ 48 Bromofluorobenzene	30	28	92.68	60-150

Metals Data

NYTEST ENVIRONMENTAL INC.

INORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

296-FB

Lab Name: NYTEST_ENV_INC. Contract: 9421415

Lab Code: NYTEST Login No.: 22633 QC Report No.22633

Matrix (soil/water): WATER
Level (low/high) : LOW
Percent Solids : 0.0

Lab Sample ID: 263307
Date Received: 11/30/94

Concentration Units (ug/L or mg/kg dry weight): UG/L

CAS No.	Analyte	Concentration	C	Q	M
7439-92-1	Lead	3.0	U	N	F

CODES :
P: ICP; F : GFAA; CV: Cold Vapor; AS: Automated Spectrophotometric
Note: A "U" in the "C" (Concentration) column indicates the analyte was not detected in this sample; "B" = Sample value greater than Instrument Detection Limit, but less than reporting limit; "NR" = Not Required.

Comments:
296-FB

NYTEST ENVIRONMENTAL INC.

INORGANICS ANALYSIS DATA SHEET

SAMPLE NO.

290-1

Lab Name: NYTEST_ENV_INC. Contract: 9421415

Lab Code: NYTEST Login No.: 22633 QC Report No.22633

Matrix (soil/water): WATER
Level (low/high) : LOW
Percent Solids : 0.0

Lab Sample ID: 263312
Date Received: 11/30/94

Concentration Units (ug/L or mg/kg dry weight): UG/L

CAS No.	Analyte	Concentration	C	Q	M
7439-92-1	Lead	3.0	U	N	F

CODES :
P: ICP; F : GFAA; CV: Cold Vapor; AS: Automated Spectrophotometric
Note: A "U" in the "C" (Concentration) column indicates the analyte was not detected in this sample; "B" = Sample value greater than Instrument Detection Limit, but less than reporting limit; "NR" = Not Required.

Comments:
290-1

NYTEST ENVIRONMENTAL INC.

INITIAL AND CONTINUING CALIBRATION VERIFICATION

Lab Name: NYTEST_ENV_INC._____ Contract: 9421415___

Lab Code: NYTEST Login No.: 22633_ QC Report No.: 22633_

Initial Calibration Source: SPEX_____

Continuing Calibration Source: SPEX_____

Concentration Units: ug/L

Analyte	Initial Calibration True	Found	%R(1)	Continuing Calibration True	Found	%R(1)	Found	%R(1)	M
Lead	90.0	97.00	107.8	50.0	51.20	102.4	52.00	104.0	F

R : Analyte Not Required

NEI FORM 3 - (1/94)

NYTEST ENVIRONMENTAL INC.

INITIAL AND CONTINUING CALIBRATION VERIFICATION

Lab Name: NYTEST_ENV_INC._____ Contract: 9421415___

Lab Code: NYTEST Login No.: 22633_ QC Report No.: 22633_

Initial Calibration Source: SPEX_____

Continuing Calibration Source: SPEX_____

Concentration Units: ug/L

Analyte	Initial Calibration True	Found	%R(1)	Continuing Calibration True	Found	%R(1)	Found	%R(1)	M
ead				50.0	53.30	106.6	52.10	104.2	F

N : Analyte Not Required

NEI FORM 3 - (1/94)

NYTEST ENVIRONMENTAL INC.

INITIAL AND CONTINUING CALIBRATION VERIFICATION

Lab Name: NYTEST_ENV_INC._____ Contract: 9421415___

Lab Code: NYTEST Login No.: 22633_ QC Report No.: 22633_

Initial Calibration Source: SPEX_____

Continuing Calibration Source: SPEX_____

Concentration Units: ug/L

Analyte	Initial Calibration True	Found	%R(1)	Continuing Calibration True	Found	%R(1)	Found	%R(1)	M
Lead				50.0	53.10	106.2	52.10	104.2	F

R : Analyte Not Required

NEI FORM 3 - (1/94)

TEST ENVIRONMENTAL INC.

INITIAL AND CONTINUING CALIBRATION VERIFICATION

ab Name: NYTEST_ENV_INC._____ Contract: 9421415___

Code: NYTEST Login No.: 22633_ QC Report No.: 22633_

nitial Calibration Source: SPEX_____

tinuing Calibration Source: SPEX_____

Concentration Units: ug/L

Analyte	Initial Calibration True	Found	%R(1)	Continuing Calibration True	Found	%R(1)	Found	%R(1)	M
ead				50.0	54.00	108.0	53.50	107.0	F

: Analyte Not Required

NEI FORM 3 - (1/94)

000058

YTEST ENVIRONMENTAL INC.

INITIAL AND CONTINUING CALIBRATION VERIFICATION

ab Name: NYTEST_ENV_INC._____ Contract: 9421415___

Code: NYTEST Login No.: 22633_ QC Report No.: 22633_

nitial Calibration Source: SPEX_____

ntinuing Calibration Source: SPEX_____

Concentration Units: ug/L

Analyte	Initial Calibration True	Found	%R(1)	Continuing Calibration True	Found	%R(1)	Found	%R(1)	M
ead				50.0	52.90	105.8	52.90	105.8	F

: Analyte Not Required

NEI FORM 3 - (1/94)

000059

YTEST ENVIRONMENTAL INC.

INITIAL AND CONTINUING CALIBRATION VERIFICATION

Lab Name: NYTEST_ENV_INC._____ Contract: 9421415___

Lab Code: NYTEST Login No.: 22633_ QC Report No.: 22633_

Initial Calibration Source: SPEX_____

Continuing Calibration Source: SPEX_____

Concentration Units: ug/L

Analyte	Initial Calibration True	Found	%R(1)	Continuing Calibration True	Found	%R(1)	Found	%R(1)	M
Lead				50.0	52.70	105.4	51.90	103.8	F

R : Analyte Not Required

NYTEST ENVIRONMENTAL INC.

INITIAL AND CONTINUING CALIBRATION VERIFICATION

Lab Name: NYTEST_ENV_INC._____ Contract: 9421415___

Lab Code: NYTEST Login No.: 22633_ QC Report No.: 22633_

Initial Calibration Source: SPEX_____

Continuing Calibration Source: SPEX_____

Concentration Units: ug/L

Analyte	Initial Calibration True	Found	%R(1)	Continuing Calibration True	Found	%R(1)	Found	%R(1)	M
Lead				50.0	51.40	102.8	51.70	103.4	F

N : Analyte Not Required

NYTEST ENVIRONMENTAL INC.

INITIAL AND CONTINUING CALIBRATION VERIFICATION

ab Name: NYTEST_ENV_INC._____ Contract: 9421415___

a Code: NYTEST Login No.: 22633_ QC Report No.: 22633_

nitial Calibration Source: SPEX_____

o tinuing Calibration Source: SPEX_____

Concentration Units: ug/L

Analyte	Initial Calibration True	Found	%R(1)	Continuing Calibration True	Found	%R(1)	Found	%R(1)	M
L ad				50.0	51.60	103.2	52.20	104.4	F

I : Analyte Not Required

NEI FORM 3 - (1/94)

000062

NYTEST ENVIRONMENTAL INC.

INITIAL AND CONTINUING CALIBRATION VERIFICATION

Lab Name: NYTEST_ENV_INC._____ Contract: 9421415___

Lab Code: NYTEST Login No.: 22633_ QC Report No.: 22633_

Initial Calibration Source: SPEX_____

Continuing Calibration Source: SPEX_____

Concentration Units: ug/L

Analyte	Initial Calibration True	Found	%R(1)	Continuing Calibration True	Found	%R(1)	Found	%R(1)	M
Lead				50.0	53.00	106.0	53.90	107.8	F

N : Analyte Not Required

NYTEST ENVIRONMENTAL INC.

INITIAL AND CONTINUING CALIBRATION VERIFICATION

Lab Name: NYTEST_ENV_INC._____ Contract: 9421415___

Lab Code: NYTEST Login No.: 22633_ QC Report No.: 22633_

Initial Calibration Source: SPEX_____

Continuing Calibration Source: SPEX_____

Concentration Units: ug/L

Analyte	Initial Calibration True	Found	%R(1)	Continuing Calibration True	Found	%R(1)	Found	%R(1)	M
Lead				50.0	53.80	107.6	54.50	109.0	F

L : Analyte Not Required

NYTEST ENVIRONMENTAL INC.

INITIAL AND CONTINUING CALIBRATION VERIFICATION

Name: NYTEST_ENV_INC._____ Contract: 9421415___

Code: NYTEST Login No.: 22633_ QC Report No.: 22633_

nitial Calibration Source: SPEX_____

tinuing Calibration Source: SPEX_____

Concentration Units: ug/L

Analyte	Initial Calibration True	Found	%R(1)	Continuing Calibration True	Found	%R(1)	Found	%R(1)	M
ead	90.0	95.60	106.2	50.0	48.30	96.6	47.90	95.8	F

: Analyte Not Required

NEI FORM 3 - (1/94)

000065

YTEST ENVIRONMENTAL INC.

INITIAL AND CONTINUING CALIBRATION VERIFICATION

ab Name: NYTEST_ENV_INC._____ Contract: 9421415___

b Code: NYTEST Login No.: 22633_ QC Report No.: 22633_

nitial Calibration Source: SPEX_____

ntinuing Calibration Source: SPEX_____

Concentration Units: ug/L

Analyte	Initial Calibration True	Found	%R(1)	Continuing Calibration True	Found	%R(1)	Found	%R(1)	M
ead				50.0	51.70	103.4	50.90	101.8	F

R : Analyte Not Required

NYTEST ENVIRONMENTAL INC.

INITIAL AND CONTINUING CALIBRATION VERIFICATION

Lab Name: NYTEST_ENV_INC._____ Contract: 9421415___

Lab Code: NYTEST Login No.: 22633_ QC Report No.: 22633_

Initial Calibration Source: SPEX_____

Continuing Calibration Source: SPEX_____

Concentration Units: ug/L

Analyte	Initial Calibration True	Found	%R(1)	Continuing Calibration True	Found	%R(1)	Found	%R(1)	M
Lead				50.0	52.60	105.2	52.40	104.8	F

NR : Analyte Not Required

YTEST ENVIRONMENTAL INC.

INITIAL AND CONTINUING CALIBRATION VERIFICATION

Lab Name: NYTEST_ENV_INC._____ Contract: 9421415___

Lab Code: NYTEST Login No.: 22633_ QC Report No.: 22633_

Initial Calibration Source: SPEX_____

Continuing Calibration Source: SPEX_____

Concentration Units: ug/L

Analyte	Initial Calibration True	Found	%R(1)	Continuing Calibration True	Found	%R(1)	Found	%R(1)	M
Lead				50.0	52.00	104.0	53.00	106.0	F

N : Analyte Not Required

TEST ENVIRONMENTAL INC.

INITIAL AND CONTINUING CALIBRATION VERIFICATION

ab Name: NYTEST_ENV_INC._____ Contract: 9421415___

b Code: NYTEST Login No.: 22633_ QC Report No.: 22633_

nitial Calibration Source: SPEX_____

ntinuing Calibration Source: SPEX_____

Concentration Units: ug/L

Analyte	Initial Calibration True	Found	%R(1)	Continuing Calibration True	Found	%R(1)	Found	%R(1)	M
ead				50.0	53.90	107.8	51.70	103.4	F

K : Analyte Not Required

YTEST ENVIRONMENTAL INC.

CRDL STANDARD FOR AA AND ICP

Lab Name: NYTEST_ENV_INC._____ Contract: 9421415___

ab Code: NYTEST Login No.: 22633_ QC Report No. : 22633__

AA CRDL Standard Source: SPEX_____

CP CRDL Standard Source: _____

Concentration Units: ug/L

	CRDL Standard for AA			CRDL Standard for ICP				
					Initial			Final
Analyte	True	Found	%R	True	Found	%R	Found	%R
Lead	3.0	3.40	113.3					

000070

NEI FORM 3 - (1/94)

ILMO2.1

YTEST ENVIRONMENTAL INC.

CRDL STANDARD FOR AA AND ICP

Lab Name: NYTEST_ENV_INC._____ Contract: 9421415___

ab Code: NYTEST Login No.: 22633_ QC Report No. : 22633__

AA CRDL Standard Source: SPEX_____

CP CRDL Standard Source: _____

Concentration Units: ug/L

Analyte	CRDL Standard for AA			CRDL Standard for ICP				
	True	Found	%R	True	Initial Found	%R	Final Found	%R
Lead	3.0	3.00	100.0					

000071

NYTEST ENVIRONMENTAL INC.

ANALYTICAL AND METHOD BLANK SUMMARY

Lab Name: NYTEST_ENV_INC._____ Contract: 9421415___

Lab Code: NYTEST Login No.: 22633_ QC Report No.: 22633__

Preparation Blank Matrix (soil/water): WATER

Preparation Blank Concentration Units (ug/L or mg/kg): UG/L_

Analyte	Initial Calib. Blank (ug/L)	C	Continuing Calibration Blank (ug/L) 1	C	2	C	3	C	Prepa-ration Blank	C	M
Lead	3.0	U	3.0	U	3.0	U	3.0	U	3.000	U	F

NR = Analyte Not Requested

NEI FORM 5 - (1/94)

000072

NYTEST ENVIRONMENTAL INC.

ANALYTICAL AND METHOD BLANK SUMMARY

Lab Name: NYTEST_ENV_INC._____ Contract: 9421415___

Lab Code: NYTEST Login No.: 22633_ QC Report No.: 22633__

Preparation Blank Matrix (soil/water): WATER

Preparation Blank Concentration Units (ug/L or mg/kg): UG/L_

Analyte	Initial Calib. Blank (ug/L)	C	Continuing Calibration Blank (ug/L) 1	C	2	C	3	C	Prepa-ration Blank	C	M
Lead			3.0	U	3.0	U	3.0	U	3.000	U	F

NR = Analyte Not Requested

NEI FORM 5 - (1/94)

YTEST ENVIRONMENTAL INC.

ANALYTICAL AND METHOD BLANK SUMMARY

ab Name: NYTEST_ENV_INC._____ Contract: 9421415___

ıb Code: NYTEST Login No.: 22633_ QC Report No.: 22633__

reparation Blank Matrix (soil/water): _____

eparation Blank Concentration Units (ug/L or mg/kg): _____

Analyte	Initial Calib. Blank (ug/L)	C	Continuing Calibration Blank (ug/L) 1	C	2	C	3	C	Prepa-ration Blank	C	M
Lead			3.0	U	3.0	U	3.0	U			F

NR = Analyte Not Requested

NEI FORM 5 - (1/94)

000074

NYTEST ENVIRONMENTAL INC.

ANALYTICAL AND METHOD BLANK SUMMARY

Lab Name: NYTEST_ENV_INC._____ Contract: 9421415___

Lab Code: NYTEST Login No.: 22633_ QC Report No.: 22633__

Preparation Blank Matrix (soil/water): _____

Preparation Blank Concentration Units (ug/L or mg/kg): _____

Analyte	Initial Calib. Blank (ug/L)	C	Continuing Calibration Blank (ug/L) 1	C	2	C	3	C	Prepa-ration Blank	C	M
Lead			3.0	U	3.0	U	3.0	U			F

NR = Analyte Not Requested

NYTEST ENVIRONMENTAL INC.

ANALYTICAL AND METHOD BLANK SUMMARY

Lab Name: NYTEST_ENV_INC._____ Contract: 9421415___

Lab Code: NYTEST Login No.: 22633_ QC Report No.: 22633__

Preparation Blank Matrix (soil/water): _____

Preparation Blank Concentration Units (ug/L or mg/kg): _____

Analyte	Initial Calib. Blank (ug/L)	C	Continuing Calibration Blank (ug/L) 1	C	2	C	3	C	Prepa-ration Blank	C	M
Lead			3.0	U	3.0	U	3.0	U			F

NR = Analyte Not Requested

NEI FORM 5 - (1/94)

YTEST ENVIRONMENTAL INC.

ANALYTICAL AND METHOD BLANK SUMMARY

Lab Name: NYTEST_ENV_INC._____ Contract: 9421415___

Lab Code: NYTEST Login No.: 22633_ QC Report No.: 22633__

Preparation Blank Matrix (soil/water): _____

Preparation Blank Concentration Units (ug/L or mg/kg): _____

Analyte	Initial Calib. Blank (ug/L)	C	Continuing Calibration Blank (ug/L) 1	C	2	C	3	C	Prepa-ration Blank	C	M
Lead			3.0	U	3.0	U	3.0	U			F

NR = Analyte Not Requested

NEI FORM 5 - (1/94)

000077

YTEST ENVIRONMENTAL INC.

ANALYTICAL AND METHOD BLANK SUMMARY

Lab Name: NYTEST_ENV_INC._____ Contract: 9421415___

Lab Code: NYTEST Login No.: 22633_ QC Report No.: 22633__

Preparation Blank Matrix (soil/water): _____

Preparation Blank Concentration Units (ug/L or mg/kg): _____

Analyte	Initial Calib. Blank (ug/L)	C	Continuing Calibration Blank (ug/L) 1	C	2	C	3	C	Prepa-ration Blank	C	M
Lead			3.0	U	3.0	U					F

NR = Analyte Not Requested

NEI FORM 5 - (1/94)

000078

YTEST ENVIRONMENTAL INC.

ANALYTICAL AND METHOD BLANK SUMMARY

Lab Name: NYTEST_ENV_INC._____ Contract: 9421415___

Lab Code: NYTEST Login No.: 22633_ QC Report No.: 22633__

Preparation Blank Matrix (soil/water): _____

Preparation Blank Concentration Units (ug/L or mg/kg): _____

Analyte	Initial Calib. Blank (ug/L)	C	Continuing Calibration Blank (ug/L) 1	C	2	C	3	C	Prepa-ration Blank	C	M
Lead	3.0	U	3.0	U	3.0	U	3.0	U			F

NR = Analyte Not Requested

NEI FORM 5 - (1/94)

000079

YTEST ENVIRONMENTAL INC.

ANALYTICAL AND METHOD BLANK SUMMARY

Lab Name: NYTEST_ENV_INC._____ Contract: 9421415___

Lab Code: NYTEST Login No.: 22633_ QC Report No.: 22633__

Preparation Blank Matrix (soil/water): _____

Preparation Blank Concentration Units (ug/L or mg/kg): _____

Analyte	Initial Calib. Blank (ug/L)	C	Continuing Calibration Blank (ug/L) 1	C	2	C	3	C	Prepa-ration Blank	C	M
Lead			3.0	U	3.0	U	3.0	U			F

NR = Analyte Not Requested

NEI FORM 5 - (1/94)

000080

YTEST ENVIRONMENTAL INC.

ANALYTICAL AND METHOD BLANK SUMMARY

Lab Name: NYTEST_ENV_INC._____ Contract: 9421415___

Lab Code: NYTEST Login No.: 22633_ QC Report No.: 22633__

Preparation Blank Matrix (soil/water): _____

Preparation Blank Concentration Units (ug/L or mg/kg): _____

Analyte	Initial Calib. Blank (ug/L)	C	Continuing Calibration Blank (ug/L) 1	C	2	C	3	C	Prepa-ration Blank	C	M
Lead			3.0	U	3.0	U	3.0	U			F

NR = Analyte Not Requested

NEI FORM 5 - (1/94)

000081

TEST ENVIRONMENTAL INC.

ANALYTICAL AND METHOD BLANK SUMMARY

ab Name: NYTEST_ENV_INC._____ Contract: 9421415___

a Code: NYTEST Login No.: 22633_ QC Report No.: 22633__

reparation Blank Matrix (soil/water): _____

reparation Blank Concentration Units (ug/L or mg/kg): _____

Analyte	Initial Calib. Blank (ug/L)	C	Continuing Calibration Blank (ug/L) 1	C	2	C	3	C	Prepa-ration Blank	C	M
Lead			3.0	U							F

R = Analyte Not Requested

NYTEST ENVIRONMENTAL INC.

MATRIX SPIKE RECOVERY DATA SHEET

SAMPLE NO.

2044-3S

a Name: NYTEST_ENV_INC._____ Contract: 9421415___

ab Code: NYTEST Login No.: 22633_ QC Report No. : 22633_

a rix (soil/water): WATER_ Level (low/med): LOW___

Solids for Sample: __0.0

Concentration Units (ug/L or mg/kg dry weight): UG/L_

nalyte	Control Limit %R	Spiked Sample Result (SSR)	C	Sample Result (SR)	C	Spike Added (SA)	%R	Q	M
Lead	75-125	25.2000		3.0000	U	20.00	126.0	N	F

omments:
2044-3_____

NR : Analyte Not Required

NEI FORM 7 - (1/94)

000083

NYTEST ENVIRONMENTAL INC.

MATRIX SPIKE RECOVERY DATA SHEET

SAMPLE NO.

296-6MSD

a Name: NYTEST_ENV_INC._____ Contract: 9421415___

ab Code: NYTEST Login No.: 22633_ QC Report No. : 22633_

a rix (soil/water): WATER_ Level (low/med): LOW___

Solids for Sample: __0.0

Concentration Units (ug/L or mg/kg dry weight): UG/L_

nalyte	Control Limit %R	Spiked Sample Result (SSR)	C	Sample Result (SR)	C	Spike Added (SA)	%R	Q	M
Lead	75-125	27.2000		5.1000		20.00	110.5		F

omments:
296-6MSD

NR : Analyte Not Required

NEI FORM 7 - (1/94)

000084

NYTEST ENVIRONMENTAL INC.

DUPLICATES

SAMPLE NO.

2044-3D

Name: NYTEST_ENV_INC._____ Contract: 9421415___

Lab Code: NYTEST Login No.: 22633_ QC Report No. : 22633__

Matrix (soil/water): WATER Level (low/med): _LOW__

% Solids for Sample: __0.0 % Solids for Duplicate: ___0.0

Concentration Units (ug/L or mg/kg dry weight): UG/L_

Analyte	Control Limit	Sample (S)	C	Duplicate (D)	C	RPD	Q	M
Lead		3.0000	U	3.0000	U			F

NR : Analyte Not Requested

NEI FORM 9 (1/94)

000085

NYTEST ENVIRONMENTAL IN[illegible].

DUPLICATES

SAMPLE NO.

296-6MS

Lab Name: NYTEST_ENV_INC._____ Contract: 9421415___

Lab Code: NYTEST Login No.: 22633_ QC Report No. : 22633__

Matrix (soil/water): WATER Level (low/med): _LOW__

% Solids for Sample: __0.0 % Solids for Duplicate: ___0.0

Concentration Units (ug/L or mg/kg dry weight): UG/L_

Analyte	Control Limit	Sample (S)	C	Duplicate (D)	C	RPD	Q	M
Lead	5.0	5.1000		4.7000	B	8.2		F

NR : Analyte Not Requested

NYTEST ENVIRONMENTAL INC.

LABORATORY CONTROL SAMPLE

Lab Name: NYTEST_ENV_INC._____ Contract: 9421415___

Lab Code: NYTEST Login No.: 22633_ QC Report No.: 22633_

Solid LCS Source: _____

Aqueous LCS Source: SPEX_____

Analyte	Aqueous (ug/L) True	Found	%R	Solid (mg/kg) True	Found	C	Limits		%R
Lead	90.0	88.40	98.2						

NYTEST ENVIRONMENTAL INC.

LABORATORY CONTROL SAMPLE

Lab Name: NYTEST_ENV_INC._____ Contract: 9421415___

Lab Code: NYTEST Login No.: 22633_ QC Report No.: 22633_

Solid LCS Source: _____

Aqueous LCS Source: SPEX_____

Analyte	Aqueous (ug/L) True	Found	%R	Solid (mg/kg) True	Found	C	Limits		%R
Lead	90.0	79.80	88.7						

000088

NYTEST ENVIRONMENTAL INC.

Instrument Detection Limits (Quarterly)

Lab Name: NYTEST_ENV_INC._____ Contract: 9421415___

Lab Code: NYTEST Login No.: 22633_ QC Report No.: 22633__

ICP ID Number: _____ Date in Effect: 10/01/94

Flame AA ID Number : _____

Furnace AA ID Number : 3030#1_____

Analyte	Wave-length (nm)	Back-ground	CRDL (ug/L)	IDL (ug/L)	M
Lead	283.30	BZ	5	3.0	F

Comments:

NEI FORM 13 - (1/94)