

**DRAFT**  
**Groundwater Sampling  
and Analysis Report**  
**2<sup>nd</sup> Quarter 2009**

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**Camp Bonneville Military Reservation  
23201 Northeast Pluss Road,  
Vancouver, WA 98682**

**Prepared For:  
Washington State  
Department of Ecology**

**Prepared By:  
Bonneville Conservation,  
Restoration & Renewal Team**

**September 2009**





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## 1.0 INTRODUCTION

This document presents the results of the Groundwater Sampling and Analysis, 2<sup>nd</sup> Quarter 2009 as part of Remedial Action Unit 2C (RAU 2C) at the former Camp Bonneville Military Reservation (CBMR) in Clark County, Washington (**Figures 1.1 and 1.2**). This report has been prepared for and is submitted by the Bonneville Conservation Restoration and Renewal Team, LLC (BCRRT), the current owner of the CBMR. The report is based on previously approved Draft Supplemental Groundwater Remedial Investigation Work Plan (GWP; BCRRT 2006, revised 2007), previous remedial investigations, previously submitted groundwater monitoring reports for Landfill 4/ Demolition Area 1 (LF4/DA1) and Site-Wide Groundwater Sentinel wells, and direction given by Washington Department of Ecology (Ecology). This submittal is part of an ongoing dialogue between the Ecology and the BCRRT regarding the applicable requirements of the Prospective Purchaser Consent Decree (PPCD; Ecology, 2006) as it relates to the RAU 2C, which addresses groundwater quality concerns.

This quarterly report includes:

- Section 1 - Introduction
- Section 2 – Site Background,
- Section 3 - Groundwater Sampling,
- Section 4 - Data management and data quality assessment (DQA),
- Section 5 - Groundwater monitoring results,
- Section 6 - Recent trends in groundwater quality,
- Section 7 - References.

This report is submitted in support of the remedial RI/FS and the specifications of regulations promulgated under the Washington State Model Toxics Control Act (MTCA) as set forth in Title 173-340 of the Washington Administrative Code (WAC) Section 350 – Remedial Investigation and Feasibility Study [WAC 173-340-350], the requirements of WAC 173-340-360 concerning the evaluation of cleanup action alternatives.

### 1.1 Purpose and Scope

The purpose of this groundwater monitoring for RAU 2C is to document and present:

- Address groundwater contamination arising from anthropogenic activities and determine the nature and extent of groundwater contamination at CBMR;
- Collect sufficient data to meet the Model Toxics Control Act (MTCA) requirements for site characterization (WAC 173-340) and other applicable groundwater monitoring guidelines published by Ecology (Ecology 1995, 2001, 2004, 2005a, 2005b and Cruz 2005).
- Generate data necessary to evaluate remedy performance; and,
- Provide an early warning should any contamination pose a threat to water supply wells or surface waters on or off CBMR.



## 1.2 General Site Information

This section contains the following general facility information:

Project title: Remedial Investigation/Feasibility Study RI/FS for Site-Wide Groundwater - Remedial Action Unit 2C

Project coordinator: Name: Michael Gage  
Address: Bonneville Conservation Restoration and Renewal Team, LLC  
23201 Northeast Pluss Road  
Vancouver, WA 98682

Phone number: (360)566-6990

Facility location: LF4/DA1 is within the boundaries of Camp Bonneville which is located in southwestern Washington; approximately 5 miles east of the Vancouver City limits in Clark County (see **Figure 1.1**).

Dimensions of facility: Camp Bonneville consists of approximately 3,840 acres. The LF4/DA1 consists of an approximately 1.82 acres located about 1,800 feet north of the Central Valley Floor and 2.5 miles northeast of where Lacamas Creek exits Camp Bonneville.

Present owner and operator: Camp Bonneville are owned and operated by the by BCRRT, LLC.

Chronological listing of past owners and operators and operational history: Since the early 1900's, the Department of the Army has owned and operated the Camp Bonneville site. In October 2006, the Army transferred ownership of the property to the County, which subsequently transferred the land to the BCRRT. BCRRT will hold the deed of the property during investigation and clean-up activities at the site. After the property is cleaned to Ecology standards the BCCRRT will transfer the property back to the County

## 1.3 Site Location and Current Land Use

The 3,840-acre Camp Bonneville site is located northeast of Vancouver, Washington, in the southeastern region of Clark County (**Figure 1.1**). The property is approximately five miles from Vancouver, Washington and approximately seven miles north of the Columbia River. Camp Bonneville is located along the western foothills of the Cascade Mountain Range, with Camp Hill and Little Elkhorn Mountain to the northwest, Munsell Hill to the west, and Little Baldy Mountain to the south.

Vehicular access to Camp Bonneville is restricted to a single entrance. The entrance is located on SE 232nd Ave. and enters the site from the west at the Camp Killpack cantonment. The entrance is gated and monitored. Most recently, the facility had been used for weekend and summer training by the U.S. Army Reserve and Navy Reserve units from Southern Washington and Northern Oregon and was a sub-installation of Fort Lewis. Other Reserve and National Guard components, as well as the Federal Bureau of Investigation (FBI) and local law enforcement units, have also used the site. Operations at the facility seized in 1995 when CBMR was selected for closure under the 1995 Base Realignment and Closure (BRAC) process. From 1995 through 2008, the FBI and other law enforcement agencies continued to use the firing range known as the



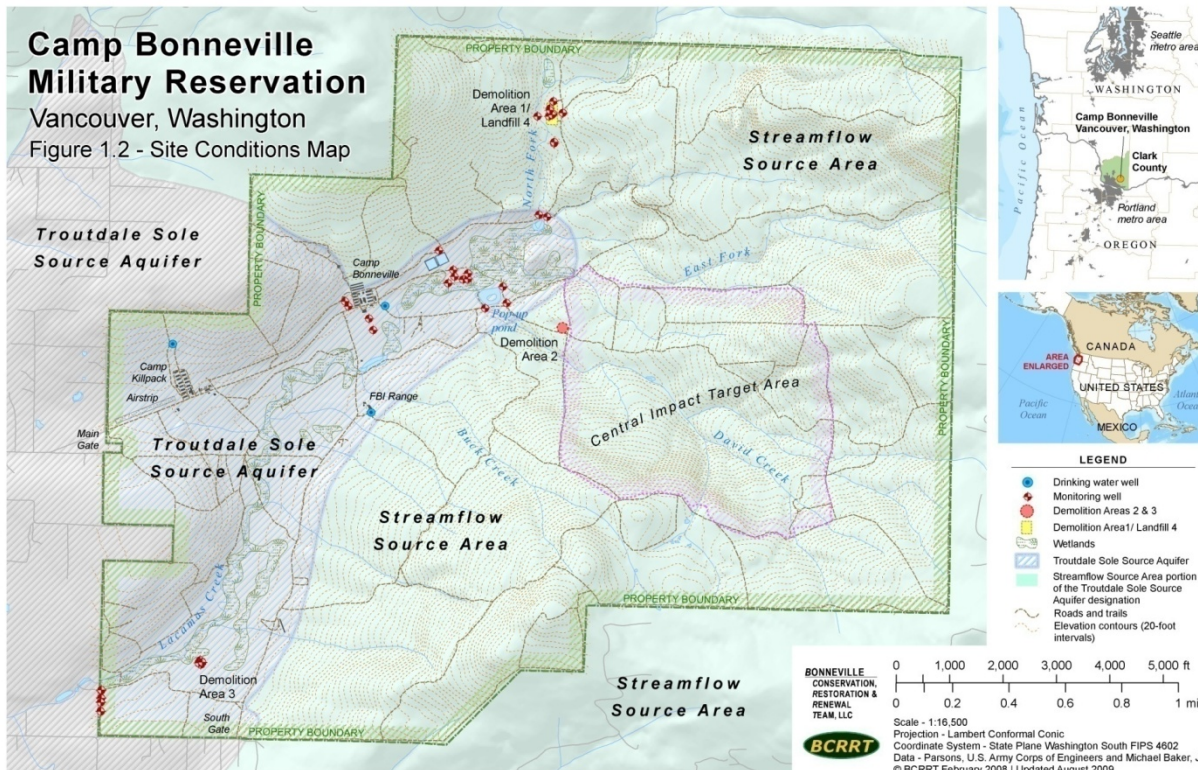
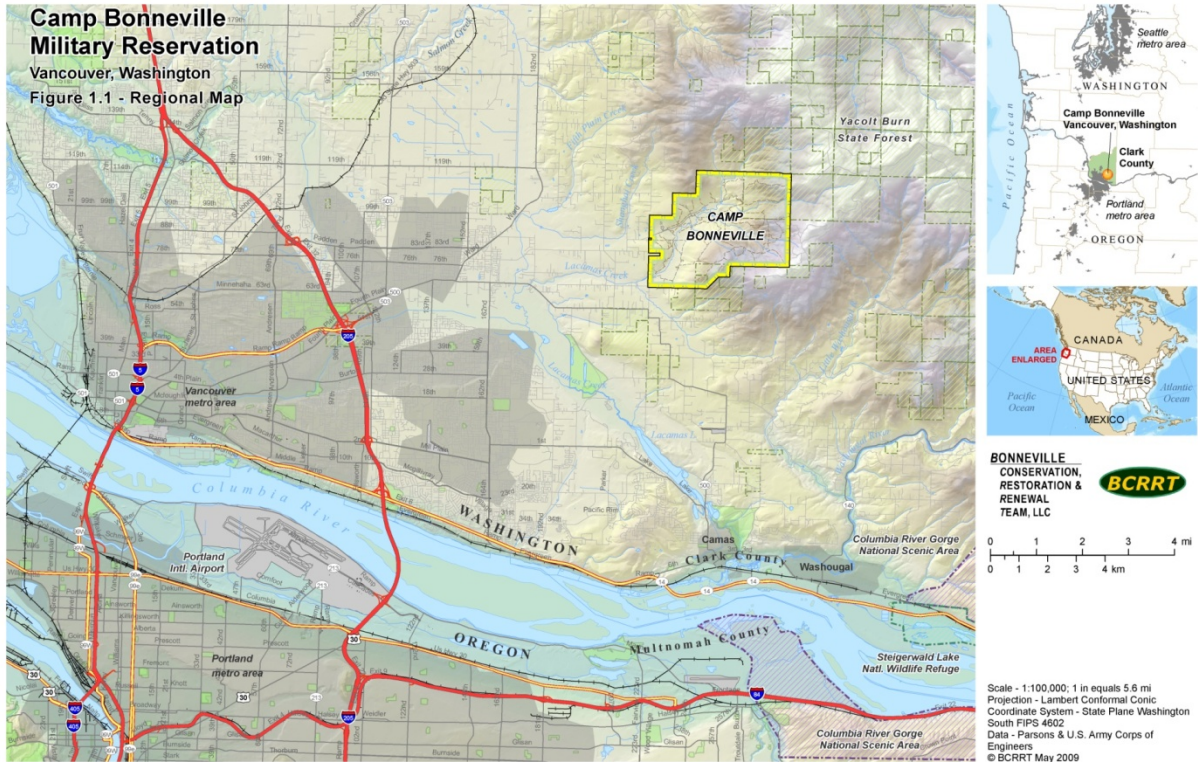


FBI Range. Currently, there is no military or law enforcement use of Camp Bonneville (grants for use of the site were cancelled beginning in November 1996). Camp Bonneville is mostly undeveloped, forested hillsides and creek side drainages. Former military barracks and training facilities are concentrated at Camp Killpack and Camp Bonneville cantonment areas. Other developed areas include firing ranges, a paved two-lane road connecting the main gate with the two cantonment areas, and a network of unpaved roads.

Between 1910 and 1995, the Army used Camp Bonneville for live fire of small arms, assault weapons, artillery, and field and air defense artillery. In the early 1950s, the Department of Defense arranged to lease an additional 840 acres from the State of Washington to expand training possibilities off of the post.

Since the Camp was officially closed, investigations have been conducted by the Army and its consultants in order to characterize the nature and extent of contamination at the site and to develop a plan for potentially transferring ownership. Clark County (County) expressed interest in the site and began the process for obtaining the property by developing a Reuse Plan. The Reuse Plan developed called for the majority of Camp Bonneville to be transferred to the County for the public benefit – education, law enforcement, and parks, with no financial gain to the County. Over the intervening years, several unsuccessful attempts were made to transfer Camp Bonneville from the Army to Clark County.

In October 2006, the Army transferred ownership of the property to the County which immediately transferred the land to the BCRRT. BCRRT will hold the deed of the property during investigation and clean-up activities at the site. After the property is cleaned to Ecology standards, BCCRT will transfer the property back to the County. The County will then begin implementing the Reuse Plan.





## 2.0 SITE BACKGROUND

The overall groundwater evaluations were summarized and discussed the BCRRT, August 2009. Draft Remedial Investigation/Feasibility Study (RI/FS) for Site-Wide Groundwater Remedial Action Unit 2C (RAU 2C), Camp Bonneville, Vancouver, Washington. This Groundwater RI/FS presented the previously approved remedial investigations, no-further-action determinations, ongoing groundwater monitoring, and direction given by Ecology. The following excerpts are from this document (shown in *italics*) and provide a brief overview the Site-Wide Groundwater Sentinel Wells and Landfill 4 /Demolition Area 1 portions of CBMR:

### “3.4 Sentinel Wells”

*A series of eight groundwater monitoring wells were installed along the valley floor at the western boundary of the CBMR. Included were: shallow wells (top of the aquifer) LCMW01S, LCMW02S, LCMW03S and LCMW04S; and deep wells (bottom of the aquifer) LCMW01D, LCMW02D, LCMW03D and LCMW04D (see **Figure 3.5**). These Sentinel Wells were constructed for the express purpose of determining whether contaminants, regardless of source, were flowing off of the CBMR at this location.”*

#### “3.4.2 Groundwater Flow”

*The shallow sets of Sentinel Wells have exhibited water level measurements that indicate a water table measured at these locations that is representative of the water table of the valley floor. The surface of the shallow aquifer is sloped from the south to the bank of Lacamas Creek, showing that water recharged on the sides of the basin flows downward into the valley where it likely turns and flows westward in parallel with the creek. Some exceptions occurred in 2004 and 2005, when the well nearest Lacamas Creek, LCMW01S suggests under certain conditions, Lacamas Creek is a losing stream in this area and is recharging local groundwater. Hence, there is a small mounding affect beneath and approximate to the creek that diminishes in a relatively short distance due to the steep slope of the terrain.*

*Depending on the season, groundwater and the Creek maybe contributing to each other along the southwestern portion of the CBMR. A comparison of water levels measured in the paired shallow Sentinel wells indicates that Lacamas Creek is a losing stream in this portion of the creek during limited periods of time. Groundwater discharges to Lacamas Creek at other times such as in December 2003 and 2004 when the deeper well (LCMW01D) has a higher water level than the shallow well (LCMW01S). The further monitoring wells are from Lacamas Creek, the more frequently they display downward gradients indicative of active recharge zones. The exceptions occur in wetter months (April and December) when some wells display upward gradients indicative of water flowing from the basin walls and surging upwards as it approaches the valley floor. This may indicate a finite thickness for water flow that may create surface seeps during wet weather.*

*In summary, groundwater flows from the sides of the basin down into the valley floor just as surface runoff would flow. Using the water level data from the OB/OD Area 3 wells in conjunction within the Sentinel Wells it is clear that groundwater flows from the east to*

*the west, draining the sub-surface basin just as Lacamas Creek drains the topographic basin. Moreover, comparison of water levels from all 27 monitoring wells installed at CBMR confirm the conceptual model that groundwater flow is analogous to the flow of surface water in Lacamas Creek and its tributaries. Water flows down from the basin walls and out the valley, leaving the site at the point where Lacamas Creek crosses the site boundary.”*

### **“3.1.2 Data Analysis**

*Groundwater samples were collected from the Sentinel Wells -- four two-well pairs (for a total of eight wells) located near the western boundary of Camp Bonneville where Lacamas Creek exits the site to the west (Site-Wide Groundwater Sentinel Wells, labeled as LC wells). Recent sampling results are summarized as follows:*

- With the use of dedicated pumps and low flow purging/sampling techniques (which are designed to obtain water samples with lower turbidity), the reported total and dissolved metals concentrations have decreased significantly. All of the total and dissolved metals detections in groundwater from these wells are were below MTCA Method A and B regulatory screening levels with the exception of results for beryllium, which have consistently been reported as estimated values (i.e., groundwater sample concentrations of beryllium detected above the Method Detection Limit (MDL) but below the Minimum Reporting Limit (MRL) of the laboratory analytical equipment and whose accuracy is limited).*
- Petroleum hydrocarbons have not been detected in any of the Sentinel Wells throughout the monitoring period of over seven years except for an isolated detection of diesel range petroleum hydrocarbons in LCMW02DW at 0.15 mg/L in January 2006.*
- Perchlorate and explosive constituents have never been detected at any of the Sentinel Wells during seven years of groundwater monitoring.*

*The groundwater samples collected and analyzed to date display no reliable evidence of any contaminant in site groundwater being present at or near the site boundary or leaving the CBMR. All explosive compounds and total petroleum hydrocarbons were below the limits of detection.*

*Perchlorate had been detected near or at the MRL of 1 ug/L it an individual well on occasion (2004, 2005, 2008 and 2009). However, these data were found to result from laboratory errors (i.e. false positives) based on one or more of the following factors: non-detectable duplicate sample, rejection by third-party independent validator, laboratory process audit, or confirmation sampling.*

*Naturally occurring levels of inorganic elements were observed in the parts per billion ranges in both shallow and deep wells. The highest observed concentrations are generally associated with groundwater samples that had a high level of suspended solids in them. The comparable filtered sample for many of the metals displayed lower concentrations, indicating the elevated reading was the result of the acid preservative*

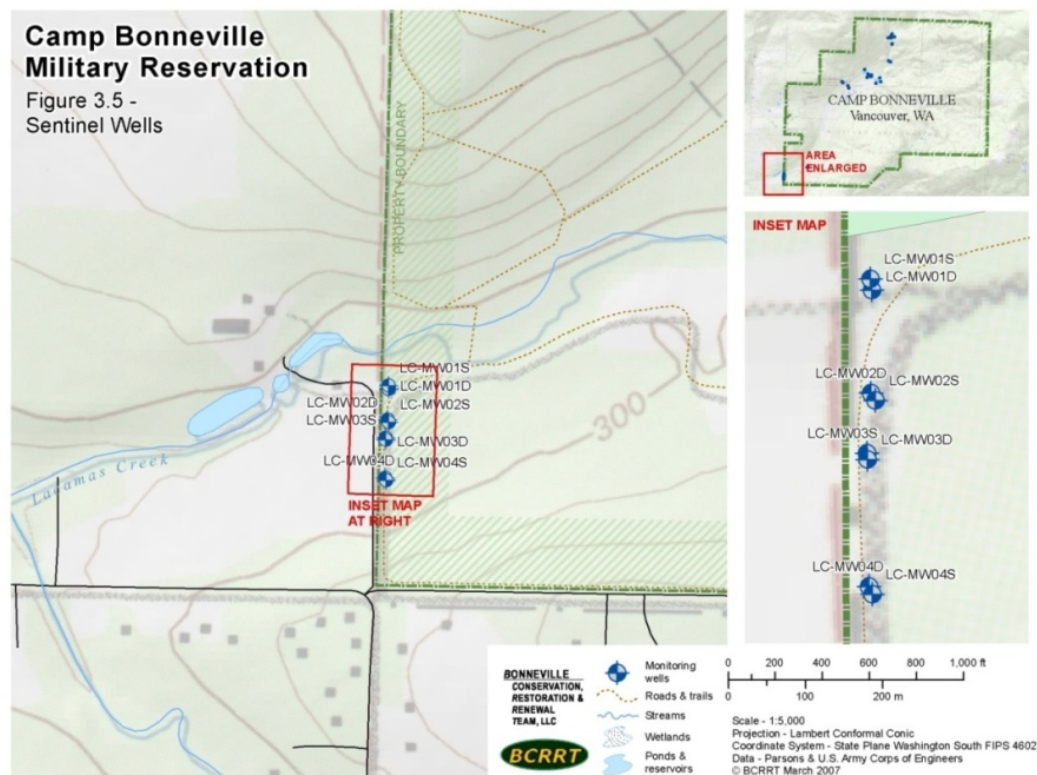
*dissolving materials from the soil particles suspended in the sample. In any event, none of the metal concentrations (total or dissolved) from Sentinel Wells exceeded MTCA Method A or Method B criteria.”*

### “3.4.3 Status Summary

*Given the extensive groundwater monitoring that has been conducted at the Sentinel Wells (located across the Lacamas Creek valley) and the lack of contaminant findings over this extended period of time, Ecology adjusted the groundwater monitoring:*

- *There are no chemicals of concern in groundwater at the Sentinel wells. No explosive, propellants, petroleum hydrocarbons, metals, SVOCs, or VOCs (except common laboratory contaminants) have been detected at levels of concern. .*
- *Four of the wells are cross-gradient and higher in elevation than the nested wells MW01S, MW01D, MW02S, and MW02D. If any groundwater contaminant plume migrates via a cross-gradient pathway (note that none have ever been detected at the CBMR boundary or the upgradient area near OB/OD Area 3) any potential contamination would be detected in these wells closer to Lacamas Creek rather than the nested wells MW03S, MW03D, MW04S and MW04D.*
- *Chemical analyses for VOCs, SVOCs, petroleum hydrocarbons (gas, diesel and oil range), and total/dissolved metals in groundwater samples from the Sentinel wells has been stopped as these analytes have never been detected in any of the sampling events conducted over more than six years..*

*While no constituents of concern have been detected, the Sentinel Wells will continue to be monitored per the Long-Term Monitoring Obligations in the PPCD.”*





### **“3.2 Landfill 4/Demolition Area 1**

*Landfill 4/Demolition Area 1 (LF4/DA1) is located about 5,000 feet northeast of the Camp Bonneville Cantonment; the landfill underlies Demolition Area 1 (see **Figure 1.2**). Vancouver Barracks reportedly used the site for the disposal of building demolition debris during the mid-1960's. According to the Environmental Baseline Survey (EBS, Woodward Clyde 1999), the former CBMR facility manager reported that firearms and ammunition were also disposed at this location. Identified potential contaminants at the CBMR included building demolition debris, explosive and propellant residue, and debris from historic onsite ordnance demolition activities, total petroleum hydrocarbons (TPH), semi-volatile organic compounds (SVOCs), volatile organic compounds (VOCs), organochlorine insecticides and herbicides, and metals. According to the Supplemental Archive Search Report (SASR, URS 1999), historic activities at CBMR included training and disposal of unserviceable ammunition. In addition, the site had been used by a number of groups and agencies, including the Army, Portland Air National Guard (PANG), local Fire Departments, and law enforcement for training and disposal operations. For example, the Bureau of Alcohol Tobacco and Firearms brought explosive and firearms to this location for disposal by open detonation.”*

#### **“3.2.1 Previous Investigations Summary**

*Site investigations (SI) were conducted to evaluate the potential for contamination resulting from past uses of the LF4/DA1. The investigation was directed primarily at evaluating potential environmental impacts from waste disposal within the landfill, but also took into account potential impacts from activities related to use of the site as an OB/OD ordnance demolition area. The primary objectives of the investigation were to evaluate whether the site poses a potential risk to human health or the environment, and to provide recommendations for additional actions (where appropriate), either for site remediation or to conduct additional investigations to better evaluate the need for and extent of remediation. The LF4/DA1 SI consisted of UXO avoidance, geophysical surveying, surface soil sampling, drilling and subsurface soil sampling, monitoring well installation and development, and groundwater sampling (URS 2003).*

*An area of buried debris disposal was identified to be approximately 120 by 200 feet during the geophysical survey. Other than a ground-penetrating radar (GPR) survey, no other types of testing were performed to delineate the actual presence of chemical constituents. The depth of the landfill material could not be determined through the use of geophysics; based on GPR profiles, it appeared to extend more than 11 feet bgs. During January 1999, groundwater was encountered at a depth of 10.4 feet below ground surface in monitoring well L4MW01 (upgradient well) and at 18.8 feet bgs in well L4MW02 (downgradient well). The report suggested that some of the landfill material could be in contact with groundwater, at least seasonally (URS 2003).*

*During the 2003 SI, the only constituents detected in soil (both near-surface and subsurface) at concentrations exceeding a MTCA Method B criteria were barium, copper, and possibly chromium. Total chromium was analyzed; however, the lowest screening criterion (which was exceeded) is based on chromium+6 (VI). Elevated barium and copper concentrations were detected in both upgradient and downgradient soil-boring samples. Arsenic, beryllium, and nickel were detected in soil samples at concentrations above MTCA Method B criteria for groundwater protections but below*

*background levels for Clark County. Low levels of one or more SVOCs, insecticides and herbicides, and VOCs were detected in some samples; however, concentrations of these constituents did not exceed the screening criteria, and several were suspected to be laboratory contaminants. Two surface soil samples were collected at RAU 2C, and 15 deeper soil samples were obtained from five borings drilled outside of the landfill area (URS 2003).*

*Initially, the only constituent detected in groundwater at a concentration exceeding a screening level was Royal Demolition Explosive (RDX, aka, cyclotrimethylene-trinitramine or hexahydro-1,3,5-trinitro-1,3,5-triazine). This compound was detected only in the presumed downgradient well (L4MW02). RDX may be associated with surface or near-surface ordnance demolition activities, rather than with deeper buried building demolition debris. Octahydro-1,3,5,7-tetranitro-1,3,5,7-tetrazocine (HMX) was also detected in L4MW02 (A), but at a concentration below the screening criteria. Low levels of three VOCs were detected in one or both wells; however, the concentrations detected were below the screening criteria. Subsequently, groundwater samples from the immediate area were found to contain perchlorate. Additional wells were installed to determine both the probable direction in which the extant plume was moving and the extent of the plume. The final two wells were installed in June of 2004 along the inferred path of migration. A total of 11 wells remain for monitoring purposes at LF4/DA1: L4MW01A; L4M201B; L4MW02A; L4MW02B; L4MW03A; L4MW03B; L4MW04A; L4MW05A; L4MW07B; L4M217; L4MW18. The A wells are installed and screened in the first water bearing strata encountered, while B wells are completed in the weathered bedrock. Wells L4MW17 and L4MW18 are relatively shallow wells completed just into the bedrock in a location where the alluvium is but a few feet thick (URS 2003).*

*Due to its relatively small size, location, explosive material hazards and potential impact to groundwater, the Army agreed to excavate and dispose of material in RAU 2C. In June 2004, a source removal action was initiated at LF4/DA1 for which the objective was to remove all fill materials and all soil contaminated above action levels. During the removal action, no construction debris was observed. In addition to materials associated with the surface ordnance disposal activities, three pits were discovered that had apparently been used for burning fireworks. Based on site observations, it appears that excess fireworks were placed in the pits and soaked with diesel oil prior to ignition. Combustion does not appear to have been complete and intact fireworks were recovered during the removal action. The pits were dug well into the heavy clay soil and one pit was completed into the saturated zone. The placement of fireworks in the saturated zone explains the observed contamination of groundwater with perchlorates in an environment that otherwise would not have significant infiltration. Following completion of the IRA excavation in September 2004 (see **Figure 3.17**), confirmatory samples indicated residual impacted soils at >20 feet below ground surface (bgs) for perchlorate, one location at > 8 feet bgs for RDX, and one location for HMX at > 5 feet bgs were present (see **Figure 3.17** and Tetra Tech 2006). Based upon the final confirmatory samples taken in December 2004, Ecology approved the completion of the soil excavation.*

*The excavation was filled to within three feet of the surface in January 2005 with clean and/or screened soils; however, the excavation was not completely filled in and graded until June 2005. Therefore, there was almost a year-long period (June 2004 to June 2005) when the excavation remained open. The open pit would have induced increased*

*recharge through the ponding of rainfall, which would have infiltrated into the fill and surrounding subsurface soils. This would have resulted in the mounding of groundwater beneath the excavation for one year and in the subsequent mobilization of many dissolved constituents that had not been mobile before.*

*During the Interim Removal Action in 2004/2005, the likely source of contaminants was found to be a series of pits dug to the water level for the disposal of fireworks (believed to have occurred in the 1991 timeframe). While most of the perchlorate and RDX source material was removed, perchlorate and RDX remain in the subsurface*

*Changes to the local hydrology occurred during and after the IRA from excavations that were 1) advanced into more permeable zones, 2) were left open and/or filled with granular material for up to a year before final backfilling and placement of topsoil/cover material, and 3) regraded the site which changed the surface runoff/recharge.*

*BCRRT subsequently conducted three soil borings and collected groundwater samples using a Geoprobe® within the perimeter of LF4/DA1 with the highest residual perchlorate concentrations in soil according to the Final Interim Removal Action Report (Tetra Tech, February 2006) and are discussed below. The locations for the borings were upgradient, east, and southeast of LF4-MW-2B (see **Figure 3.8**).*

*In addition to the LF4/DA1 sampling, surface water samples were collected from Ecology selected locations in the North Fork of Lacamas Creek where there is a potential to receive groundwater from LF4/DA1. Water samples were collected from the center of the water column and along the eastern bank. Sampling and perchlorate analysis of three surface water locations in the North Fork of Lacamas Creek – (1) upstream/northwest of MW-4A, (2) directly across/west of LF4-MW2A&B pair, and (3) downstream/south where the creek goes through two 90°bends and the mapped residual Troutdale bedrock pinches out (see **Figure 3.8**).*

*Soil, groundwater, and surface water samples were collected for laboratory analysis for perchlorate and geotechnical/in-situ analysis for hydraulic conductivity (see **Appendix D**):*

- Soil samples collected at various depths analyzed for perchlorate contained concentrations ranging from non-detect (fill material) to 2,100 ug/kg. Residual perchlorate was detected in the unsaturated soil. The upper zones of the borings indicated sandy silt while the lower zones indicated granular sand with some silt. The concentrations of perchlorate were consistent with the IRA confirmation samples (Tetra Tech 2006).*
- Shallow groundwater perchlorate concentrations were 420 ug/L and 760 ug/L while a deeper sample was 350 ug/L.*
- Three surface water samples and one field duplicate were collected from Lacamas Creek were all non-detectable for perchlorate.*
- Three geotechnical samples revealed the following concerning the following about*

*subsurface soils at LF4/DA1*

- 30' to 36' is a Silty Sand (SM) with a specific gravity of 2.757, a soil porosity of 66.0%, and composed of 55.7% sand, 34.1% silt and 10.2% clay,
  - 36' to 39' is a Silty Sand (SM) with a specific gravity of 2.756, a soil porosity of 64.6%, and composed of 60.3% sand, 33.4% silt and 6.3% clay, and
  - 40' to 41' is a Sandy Silt (ML) with a specific gravity of 2.769, a soil porosity of 65.7%, and composed of 49.8% sand, 38.8% Silt and 11.4% clay.
- *Hydraulic conductivity results from geotechnical samples, in-situ aquifer tests, and previous aquifer tests performed by others are included as **Appendix D.***

### **“3.2.2 Site Specific Conditions**

#### **3.2.2.1 Surface Water Hydrology**

*The North Fork of Lacamas Creek bounds the LF4/DA1 to the north, west, and southwest and flows to the south where it enters the Central Valley and joins with the East Fork. The water quality of Lacamas Creek is monitored indirectly south of the LF4/DA in two monitoring wells that straddle the North Fork of Lacamas Creek (where it enters the Central Valley) and in the Sentinel Wells located where Lacamas Creek exits Camp Bonneville.*

*According to the March 2000 Final Project Completion Report, Surface Water Investigation of Lacamas Creek, Camp Bonneville, Vancouver, Washington, (Hart Crowser, 2000), the results of water samples collected from Lacamas Creek and North Fork of Lacamas Creek indicated that LF4/DA1 has not impacted surface water quality..*

*While the LF4/DA1 area is technically within the Streamflow Source Area for the Troutdale Sole-Source Aquifer System, it is in that portion of the Streamflow Area that contributes via surface water recharge of the Central Valley sediments – no groundwater recharge is believed to occur. The groundwater in the LF4/DA1 area does not directly connect to the Central Valley but may discharge into the North Fork of Lacamas Creek, where it would undergo significant dilution based on the size of the creek's recharge basin vs the LF4/DA1 recharge area.*

#### **3.2.2.2 Geology**

***Figures 2.2 and 3.8** show that the LF4/DA1 site is isolated geologically and is surrounded by wide expanses of the impermeable andesite bedrock (Them). Essentially the LF4/DA1 is located within a geologic “island” comprised of the conglomerate (see **Section 2.4.2**) members of the Troutdale Formation (Ttfc), and there is no connection of the LF4/DA1 to the Troutdale units on the western edge of the CBMR.*

*In three dimensions, the geology of the LF4/DA1 area may be more accurately described as a Ttfc “cap” on a hillside otherwise consisting of Them. However,*

*little remains of the original appearance of either rock formation due to extreme weathering. Well logs show that both units exist as sands, silts, and clays to a depth of about 70 feet. The conglomerate member of the Ttfc has been weathered to clayey silt and the andesite bedrock to sandy clay (see **Section 2.4.2** and well logs in **Appendix B**).*

### 3.2.2.3 Hydrogeology

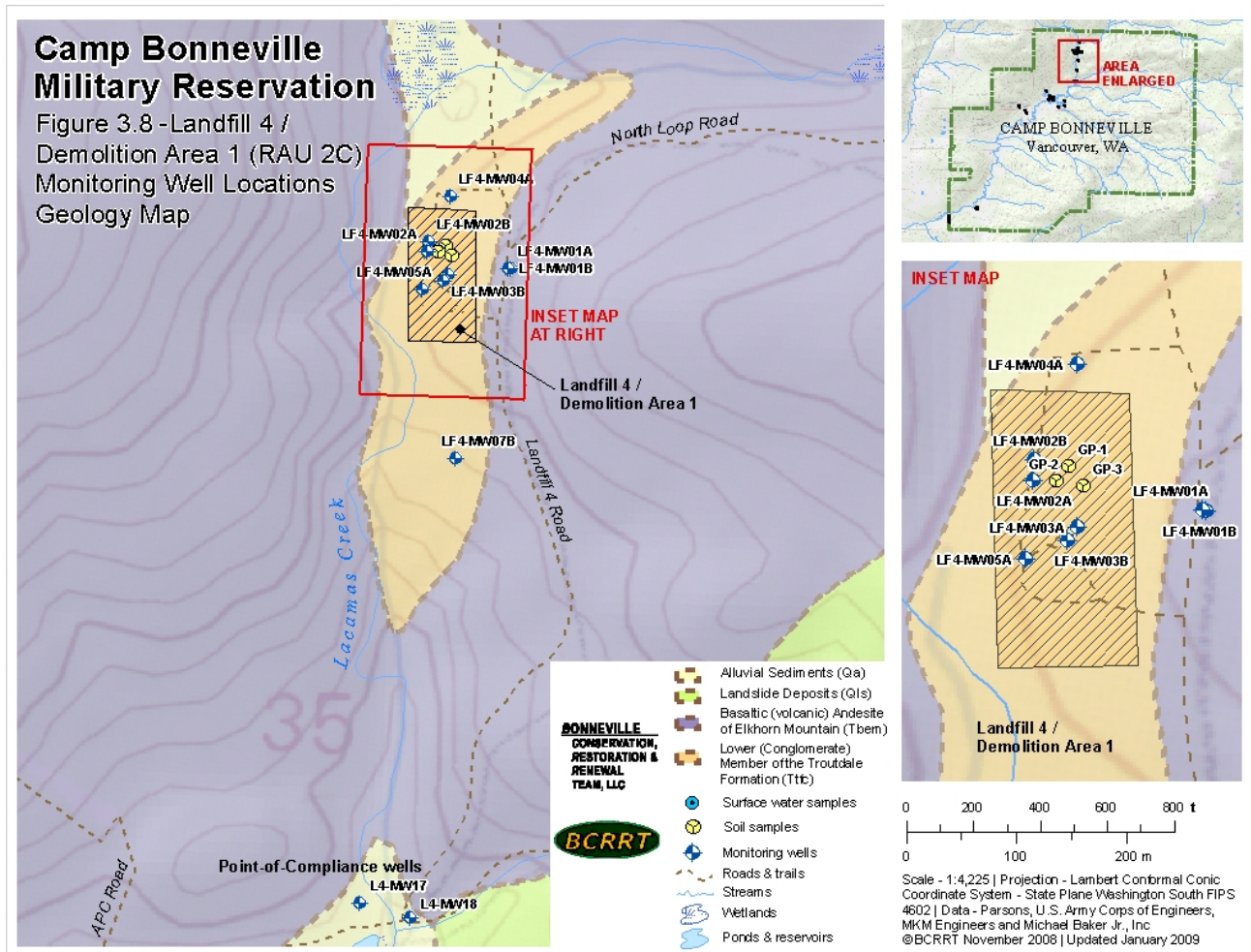
*LF4/DA1 is located in an upland area of Camp Bonneville, north of the Central Valley. **Figure 3.8** presents is a geologic map superimposed upon a topographic map and shows that there is a remnant valley fill of Ttfc (conglomerate member of the Troutdale Formation) beneath the site. However, this formation extends only to a maximum depth of about 15 feet near MW-2B and is not saturated. Groundwater occurs only within the saprolitic soil formed from the heavily weathered andesitic bedrock of the Elk Mountain basalt (Tbem)*

*A total of ten monitoring wells were installed around and near LF4/DA1 as part of the investigation of RAU 2C –LF4/DA1 (URS, 2003). Five of the wells (LF-MW-1A, 2A, 3A, 4A, 5A,) were screened in shallow soils (30 to 45 feet bgs) and four wells (LF-MW-1B, 2B, 3B, 7B) were screened in deep soils atop competent bedrock (50 to 72 feet bgs). Another shallow well (MW-6A) was installed to bedrock on the steep hillside west of the landfill and east of Lacamas Creek, but was reported as dry and has not been used. In the most recent sampling event (September 2008), the depths to water in the shallow wells ranged from 17 to 31 feet and those in the deep wells from 14 to 41 feet (see **Appendix B**).*

*At the LF4/DA1 site, all of the saturated overburden material encountered was saprolite and heavily weathered from the parent material (andesitic basalt with zeolite inclusions) into sandy silt or silty sand with white mottling. This overburden material graded into increasingly larger grain sizes with depth until competent rock was reached. There were zones of saturated sandy, silty, or clayey (angular) gravels atop the competent bedrock. Within the competent bedrock, open fractures were noted in the three rock cores (LF-MW-1B, 2B, and 3B) at LF4/DA1, most of which were reported as being oriented horizontally or nearly so (URS, 2003).*

*Such rock characteristics tend to direct groundwater flow horizontally within the overburden and the relatively shallow weathered zones of bedrock until it can move upwards in response to a hydraulic discharge point of lower pressure head (e.g., a stream). Therefore, groundwater occurrence in the bedrock is generally in the uppermost weathered/fractured zones and especially in the gravelly portions atop the competent rock.”*



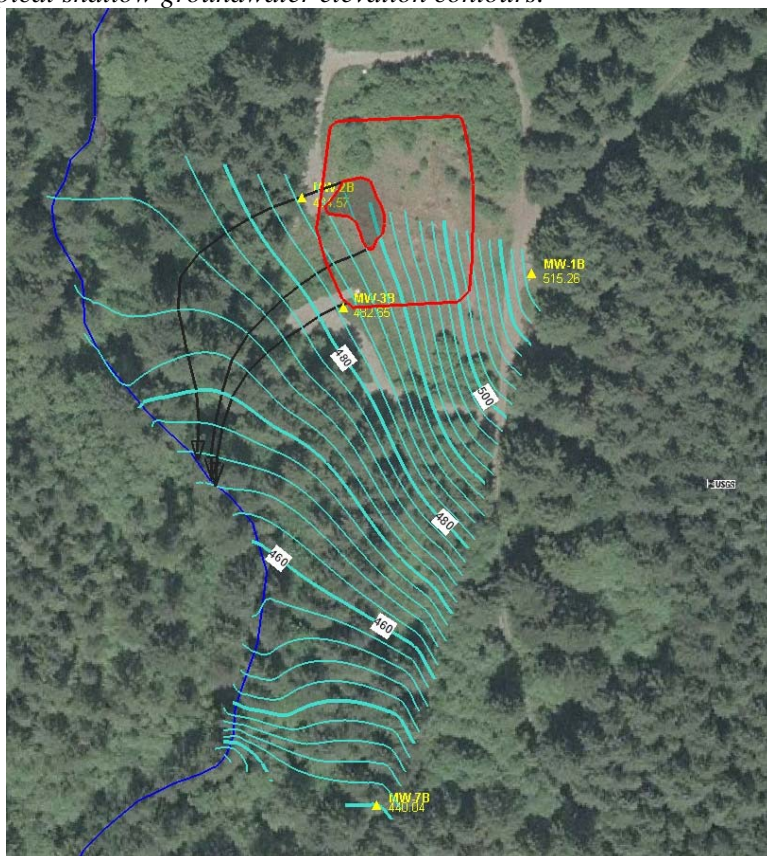


“3.2.2.4 Groundwater Flow

Groundwater elevation data from a typical sampling event (September 2008, see **Appendix B**) were mapped for groundwater flow direction analysis in the following figures. **Figure 3.9** is typical of the historical groundwater flow pattern in the shallow zone and it indicates that shallow groundwater flows west from LF-MW-1A until a groundwater divide is reached, where the flow separates toward either north (toward LF-MW-4A) or south (toward LF-MW-3A and 5A). **Figure 3.10** is typical of the deeper zone and indicates that the deeper groundwater flows west from MW-1B in a semi-radial fashion toward the North Fork of Lacamas Creek, which is consistent with the surface topography. Therefore, groundwater from the LF4/DA1 site is expected to discharge to the North Fork of Lacamas Creek within 300 feet west and/or 250 feet southwest of the LF4/DA1 site.”



*Figure 3.9 – Typical shallow groundwater elevation contours.*



*Figure 3.10– Typical deep groundwater elevation contours.*

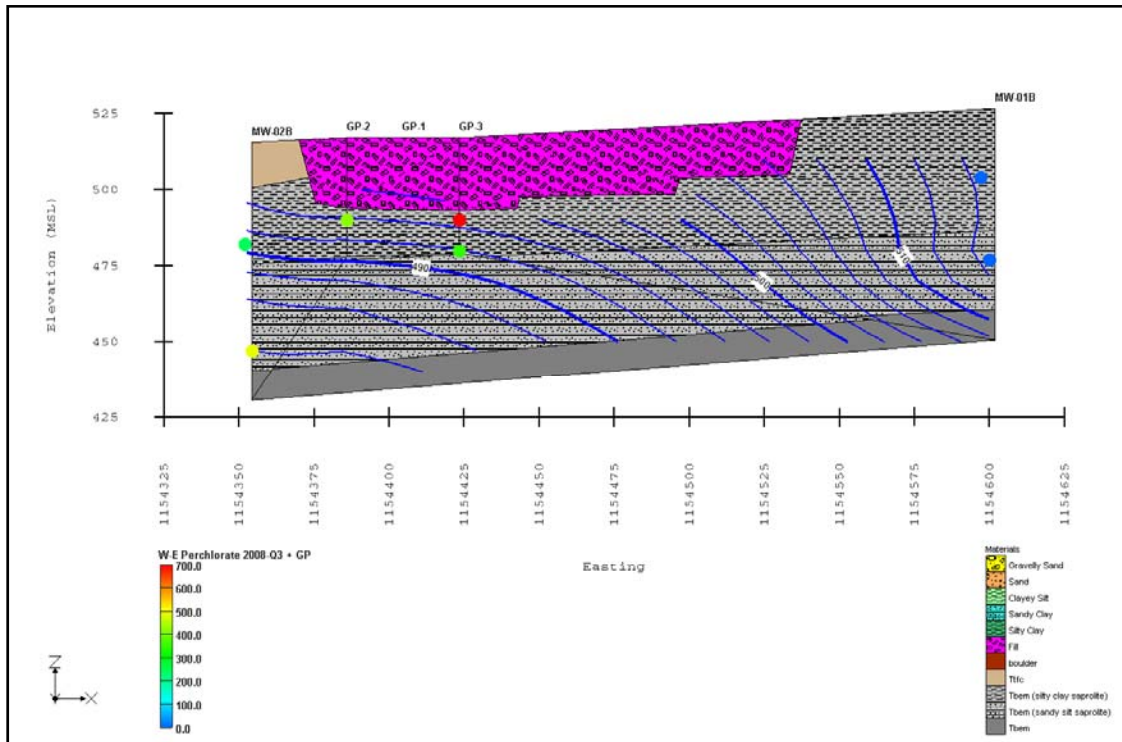


Two hydrogeologic cross-sections were also constructed using the same data set. **Figure 3.11** is a south-to-north cross-section from LF-MW-5A to 4A showing the groundwater divide in more detail, including the vertical dimension. LF-MW-2B lies just south of the groundwater divide. **Figure 3.12** shows a west-to-east cross-section from LF-MW-1A to 2B. This figure shows the change in the vertical flow component across the site: near LF-MW-1B, the flow is mostly horizontal. Toward the west, the flow becomes increasingly vertical. It appears that increased infiltration through the relatively permeable fill in the formerly excavated area is causing this alteration in the vertical flow gradients. **Figure 3.12** also may demonstrate why LF-MW-2B groundwater contains the highest perchlorate/RDX concentrations: the bedrock surface dips toward the west and, as impacted groundwater flows along the top of competent rock, it flows near well LF-MW-2B.

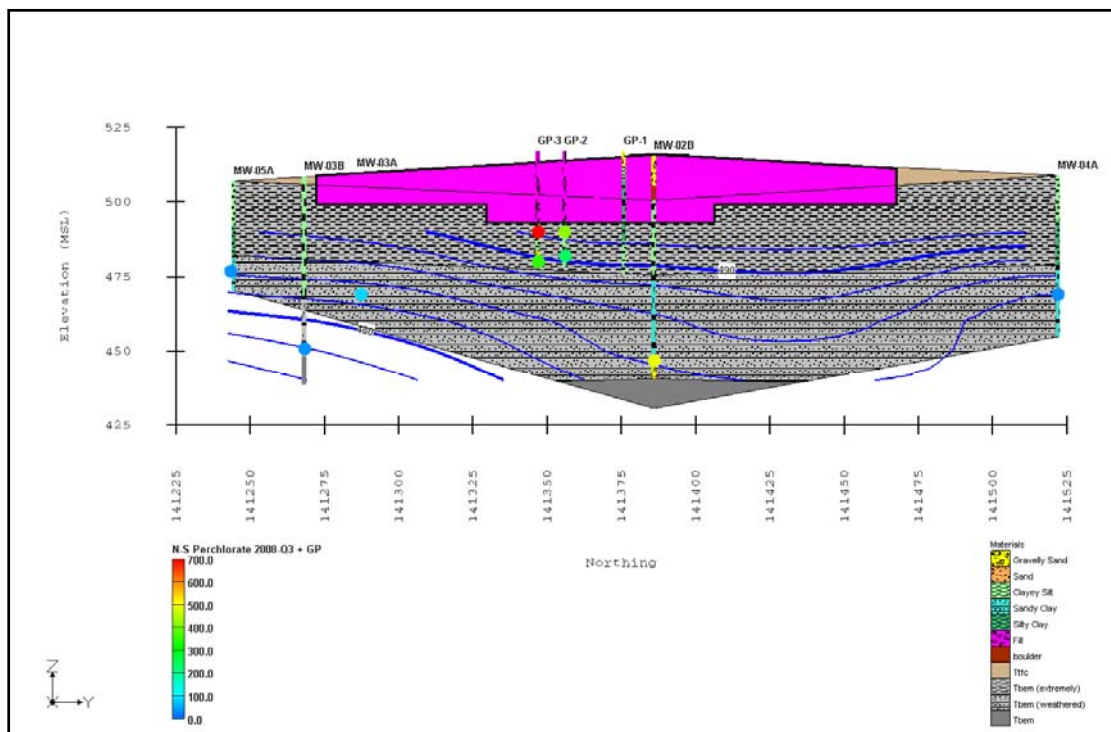
Estimates of the hydraulic conductivity of the saturated zones beneath LF4/DA1 had been made previously by using slug tests in several wells (URS 2003). However, the previous report questioned some of their own results as not corresponding to the observed stratigraphy (i.e., relatively high K values in silts and clays); therefore, the accuracy of their slug test data and/or analyses is doubtful. The raw slug test data and the corresponding digital AqTeSolv<sup>®</sup> files for those tests are not currently available for re-analysis.

A step-drawdown “yield” test had also been performed (URS 2003) in LF-MW-2B, but the data were not evaluated for hydraulic conductivity; fortunately, those data were available and were re-analyzed. The pumping rate and drawdown/recovery data were input to AqTeSolv<sup>®</sup> (version 3.5 Professional, Duffield, 2002) and analyzed using the Theis method for variable-rate pumping tests in an unconfined aquifer (a modification of Theis, 1935). The drawdown data in the pumping well were assumed to be greater than that in the formation due to friction in the well. The result was a good graphical match with the late recovery data, which are not affected by friction losses during the pumping phase. The result was a transmissivity value (T) of 9.0 ft<sup>2</sup>/d, which, when divided by the aquifer saturated thickness (b) of 44 feet, yields a hydraulic conductivity value (K) of 0.204 ft/d (7.2e-5 cm/sec).

This K value will be used along with other values derived from the Supplemental Data Collection slug tests conducted recently to obtain a representative K value for fate calculations. GP-3 was slug-tested three times each at two different depths (29 feet and 38 feet bgs). The mean K value for the 29-foot depth interval was 0.37 ft/day and the mean K of the 38-foot depth interval was 0.233 ft/day (BCRRT August 2009).



**Figure 3.11**– South-to-North Cross-Section at LF4/DA1 - Perchlorate Concentrations (ug/L) in Groundwater and Elevation Contours (Interval = 2 ft.)



**Figure 3.12** – West-to-East Cross-Section at LF4/DA1 - Perchlorate Concentrations (ug/L) in Groundwater and Elevation Contours (Interval = 2 ft.)



3.7.3.2 LF4/DA1 Wells – Characteristics Derived from Groundwater Trends

The seasonal changes at L4-MW- 2A, 2B, and 3A, mounding at 1A, and the continued presence of mobile contaminants all support the observed hydrogeology conditions that the LF4/DA1 is acting in a manner analogous to a “bathtub”. The “bathtub” is where water entering the area via infiltration becomes trapped because of the lower permeability “walls” and competent bedrock “floor” with no fast way out. The accumulated water causes mounding which slowly seeps out, taking dissolved contaminants with it.

- *The observed seasonal variations in the shallow wells reflect dilution rather than migration out of the area. The wet/Winter conditions dilute the LF4/DA1 concentrations as fresh water recharges into the area; then as groundwater adsorbs, evaporates, and seeps out, the concentrations increase to peak levels during the dry/Summer conditions.*
- *The recent 4th Quarter 2008 spike in perchlorate concentration at the upgradient L4-MW-1A is directly attributable to the heavy precipitation event prior to and during the sampling event. The influx of water into the system created a mounding effect that forced impacted groundwater upgradient. The creation of this mounding reflects the inability of groundwater to easily migrate out of the local system.*
- *The generally stable presence of very mobile constituents such as perchlorate, chlorinated solvents, and RDX in an environment that experiences significant rainfall, lacks an overlying confining layer, and has a long timeline since the placement/release of source material all indicate that the local conditions have to be retarding the migration of contaminants out of the LF4/DA1 area.*

**Table 3-2  
 Perchlorate and RDX Maximum and Latest Detections by Well**

COCs	Perchlorate <sup>(1)</sup>		RDX <sup>(1)</sup>	
	Maximum	June '09	Maximum	June '09
L4MW01A	<b>36</b>	2.9	0.49	0.13
L4MW01B	1.2	1.1	<b>0.92</b>	ND
L4MW02A	<b>280</b>	<b>195</b>	<b>40</b>	<b>19</b>
L4MW02B	<b>530</b>	<b>431</b>	<b>120</b>	<b>84</b>
L4MW03A	<b>120</b>	<b>83</b>	<b>13</b>	<b>9.4</b>
L4MW03B	<b>53</b>	<b>42</b>	<b>6.1</b>	<b>4.1</b>
L4MW04A	<b>40</b>	<b>34</b>	<b>2.8</b>	<b>2.8</b>
L4MW05A	<b>64</b>	<b>36</b>	<b>5.2</b>	<b>4.1</b>
L4MW07B	3 <sup>(2)</sup>	2.3	ND	ND
L4MW017	ND	ND	ND	ND
L4MW018	ND	ND	ND	ND

<sup>(1)</sup> **Bold** values exceeded MTCA Method A or B Cleanup Levels.

<sup>(2)</sup> Resampled value to correct a field cross-contamination result of 20 ppb that has not been repeated in the subsequent 8 quarterly sampling events.





The Section 4.0 Conceptual Site Model/Risk Assessment concluded that:

**“4.3 Summary**

*In order for there to be a risk, a complete environmental pathway by which chemicals may be transported to human or ecological receptors must exist. The only potentially complete pathways for constituents in groundwater moving away from LF4/DA1 are:*

- 1) *Future groundwater receptor, in the event that existing or new on-site wells (deep/bedrock wells) are used as a potable water source in the Central Valley.*
- 2) *Surface water for human health receptors (recreational users/fishers) via Lacamas Creek.*
- 3) *Surface water for ecological receptors in Lacamas Creek.*

The Section 5.0 Cleanup Standards proposed the following point of compliance and cleanup levels:

**“Table 5-2. Summary of Points of Compliance and Applicable Cleanup Levels.**

	<b>Landfill 4/Demolition Area 1</b>			<b>Site-Wide Groundwater</b>
<i>Point of Compliance:</i>	<i>LF4MW 1A&amp;B, LF4MW 2A&amp;B, LF4MW 3A&amp;B, LF4MW 4A, LF4MW 5A, and LF4MW 7B</i>	<i>Surface Water Monitoring Points</i>	<i>LF4MW 17 and LF4MW 18</i>	<i>LCMW 1A&amp;B, LCMW 2A&amp;B, LCMW 3A&amp;B, and LCMW 4A&amp;B</i>
<i>MTCA Cleanup Level:</i>	<i>Model/Risk Based Method C<sup>1</sup></i>	<i>Modified Method B<sup>2</sup></i>	<i>Method B</i>	<i>Method B</i>
<b>Perchlorate</b>	<i>1,300 ppb</i>	<i>600 ppb</i>	<i>11 ppb</i>	<i>11 ppb</i>
<b>RDX</b>	<i>95 ppb</i>	<i>59 ppb</i>	<i>0.8 ppb</i>	<i>0.8 ppb</i>

Notes:

<sup>1</sup> *Based upon the Section 3.7.3.5 groundwater modeling and Section 4 risk assessment criteria.*

<sup>2</sup> *Based upon the lowest cleanup level for Method B human health or ecologic receptors.”*



### 3.0 GROUNDWATER SAMPLING

Baker conducted groundwater sampling for the 2nd Quarter 2009 events at 19 existing monitoring wells at the Sentinel Wells and LF4/DA1 Wells at CBMR (see **Figures 1.2, 3.5 and 3.8**). The sampling event was conducted from June 24 to June 29, 2009. All groundwater sampling was conducted in accordance with the procedures established in the previously approved Draft Supplemental Ground Water Remedial Investigation Work Plan (groundwater workplan/GWP; BCRRT November 2006, revised September 2007). The procedures detailed in the GWP include sample collection, sample labeling, chain-of-custody, field documentation, decontamination, and investigative-derived waste (IDW) handling. All on-site activities were performed in accordance with the CBMR Site Health and Safety Plan (HASP; BCRRT October 2006, revised August 2007)

Ecology approved modifications to the GWP included:

- Installation of dedicated bladder pumps in each of the wells on March 2008 (per the letter from Baker to Ecology, dated February 13, 2008).
- Reduction in the groundwater monitoring parameters based on historic sampling results (approved in a letter by Ecology, dated March 18, 2009):
  - VOCs via USEPA Method 8260B (without tentatively identified compounds/TICs)
  - Explosives including picric acid, nitroglycerin and PETN via USEPA Method 8330
  - Perchlorate via USEPA Method 314.1
  - Field Measurements of temperature, specific conductivity, dissolved oxygen, pH, oxidation reduction potential and water levels, via calibrated field instruments.

The Sentinel Wells will also be sampled annually for priority pollutant metals, semi-volatile organic compounds (SVOCs), polynuclear aromatic compounds (PAHs), and pH.

#### 3.1 Well Purging and Sampling

Groundwater was collected via low-flow/minimal-drawdown well sampling techniques during purging and sampling. All of the wells have dedicated Solinst bladder pumps installed. The pumps are constructed with a PVC body and a Teflon bladder that is actuated via a surface controller and air supply (compressor) that are transported between the wells. The low flow purging technique is described in detail in the GWP. Low-flow sampling minimizes disturbance to the aquifer and is designed to ensure that samples collected from the wells are representative of the aquifer (without undue agitation/sediment).

At the well identified as LF4-MW-17, purging was discontinued when well was pumped dry. The well was allowed to recharge and samples were collected approximately 2-3 hours after purging began.

#### 3.2 Sample Collection

Samples were collected into the laboratory-supplied sample containers directly from the end of the dedicated discharge tubing. Appropriate sample preservative had been supplied in the containers by the laboratory. Sample containers for VOCs were filled completely to the top of the container with no headspace/bubbles, to prevent sample dilution due to volatilization.

Samples were stored in coolers with ice and maintained in secured field vehicles or field barracks prior to



daily shipment to the laboratory. Samples were maintained at 4°C during all phases of sample storage, prior to analytical testing.

Proper Chain-of-Custody (COC) documentation was maintained for all samples from the time of collection until the samples were shipped to the laboratory. COCs accompanied all of the samples and contained the following information: project number, sample numbers, number of containers, method(s) of preservation of samples, date and time of sample collection, analysis(es) requested, date and time of transportation to the laboratory, method of transportation, and any other information pertinent to the samples.

Samples were placed in appropriate shipping containers (i.e., coolers) and properly packaged to prevent damage to the samples. All samples were re-counted, the sample container labels checked against the field daily logbook and the COCs before each cooler was sealed. The completed COCs were sealed in plastic bags and placed in each cooler. Samples were picked up daily at the site by the analytical laboratory.

### **3.3 Quality Assurance/Quality Control Samples**

As specified in the Ecology approved GWP, field QA/QC samples were collected for the purpose of assessing the quality of sampling and accuracy and precision of analyses. These QA/QC samples were submitted blind to the analytical laboratory. Field duplicate samples were collected at a frequency of 1 per 10 monitoring well samples. Matrix spike/matrix spike duplicate (MS/MSD) samples were collected at a frequency of 1 per 20 monitoring well samples. Trip blanks were submitted with each cooler that contained samples for VOC analysis.

The use of dedicated pumps in all of the wells eliminated the need for equipment rinsate or field blanks.



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## 4.0 DATA MANAGEMENT/ QUALITY ASSESSMENT

All of the analytical data was received from the laboratory in an electronic data deliverable (EDD) format for importation into the CBMR databases. Laboratory and any validator qualifiers are included with the data sets. Complete copies of the laboratory data packages and validation reports are included electronically on the attached CD as **Appendix A**.

All analytical services were performed in accordance with the Ecology approved GWP. Site QA/QC procedures to ensure are fully detailed in the GWP and include:

- Internal and external audits as part of regulatory agencies laboratory monitoring;
- Following field and laboratory objectives for Precision, Completeness, Representativeness, and Comparability,
- Adhering to Sampling, Custody Procedures, and Final Evidence Files.
- Use of appropriate Field Instruments and Laboratory Instruments Calibration, Field Measurements and Laboratory Analytical Methods, field QA/QC samples, and Field and Laboratory Quality Control Checks,
- Field Performance and System Audits,
- Use of Instrument Preventative Maintenance protocols, and
- Timely Corrective Actions

BCRRT technical staff reviewed the analytical data and support documentation and determined that the data was usable. Trip blank samples submitted were non-detectable for VOCs. Field duplicate samples were within the relative percent differences (RPD) goals established in the GWP. A third party data validation was conducted for the explosives parameters analysis, in accordance with the GWP and USEPA methods.

The analytical tables include the appropriate State of Washington MTCA levels for comparison with regulatory and risk-based criteria. MTCA Method A and B Cleanup values and applicable, relevant and appropriate state and federal groundwater screening values were obtained from the Ecology CLARC database (<https://fortress.wa.gov/ecy/clarc/Reporting/CLARCReporting.aspx>). MTCA Method B-Modified and Method C Cleanup values were established in the Draft RI/FS for Site-Wide Groundwater (BCRRT August 2009) currently under review by Ecology.

### 4.1 Field Data Quality Assessment

While there are no specific data quality objectives for the measurement of field parameters (such as temperature, pH, ORP, conductivity, dissolved oxygen, and turbidity), stabilization was considered reached when three consecutive readings were within  $\pm 0.3$  for pH,  $\pm 1$  degrees C for temperature,  $\pm 10$  percent for specific conductance,  $\pm 10$ mV for ORP, and  $\pm 0.5$  mg/L for DO. If five well volumes were purged from the subject well and stabilization had not been met, sampling occurred regardless of field parameter stabilization.

### 4.2 Quality Control Sample Assessment

Trip blanks accompanied the groundwater samples for VOC analysis that were consolidated daily into one cooler and shipped to the laboratory. Trip blanks were shipped on June 24, June 25, June



26 and June 29, 2009. All four trip blanks were analyzed for VOCs and none had compounds detected above the method detection limit.

One duplicate sample was collected from each of the study areas. The duplicate samples were analyzed for the same constituents as the source sample. The RPD was calculated as the difference between the values divided by the average of the values. A significant difference between duplicate values for a few parameters indicates potential problems with the precision of specific analyses. A significant difference for many parameters indicates potential problems with the sample-collection procedures.

**Table 4-1  
 Duplicate Sample Results and Relative Percent Difference**

Site Sample ID Sample Date/Time Units	23LCMW04SW 6/24/2009 ug/L	23LCMW0460W 6/24/2009 ug/L	Relative Percent Difference	23L4MW03BW 6/26/2009 ug/L	23L4MW460W 6/26/2009 ug/L	Relative Percent Difference
1,1,1-Trichloroethane	ND	ND	--	ND	ND	--
1,1-Dichloroethane	ND	ND	--	ND	ND	--
1,1-Dichloroethene	ND	ND	--	ND	ND	--
1,3,5-Trinitrobenzene	ND	ND	--	ND	ND	--
2,4-Dinitrotoluene	ND	ND	--	ND	ND	--
2-Nitrotoluene	ND	ND	--	ND	ND	--
Dichlorodifluoromethane	ND	ND	--	ND	ND	--
HMX	ND	ND	--	ND	0.029 J	--
Perchlorate	ND	ND	--	41.9	40.5	3.34
RDX	ND	ND	--	4.1	4.1	0.00
Toluene	ND	ND	--	ND	ND	--
Trichloroethene	ND	ND	--	ND	ND	--

The Sentinel well selected for blind field duplicate analysis was LCMW-04S; the field duplicate was identified as LCMW-0460. There were no detections in the sample or the associated field duplicate located at the Sentinel Wells at Lacamas Creek; therefore no values were calculated for RPDs. Another blind field duplicate was collected from L4MW-03B; the duplicate was identified as L4MW-460B. Perchlorate was detected at 41.9 and 40.5 ug/L, respectively. The comparison of the perchlorate results indicates a RPD of 3.34% well below the 20% threshold indicated in the GWP for an acceptable difference. RDX was also detected in the LF4 sample and associated field duplicate; both results were 4.1 ug/L; there is no difference in the RPDs.

### 4.3 Laboratory Analysis Chemical Data Quality

The method blanks for the explosives data set was determined to exhibit contamination for the constituent 1,3,5-trinitrobenzene. This contamination resulted in 80% of the samples reporting a false positive results for 1,3,5 trinitrobenzene. Baker decided to have an independent data validator review all of the explosives analysis to determine whether there was any data quality issues associated with this method blank contamination. The independent validator determined that these results should be reported as a non-detect at the method detection limit; no other qualifiers were required by the validator.

During the groundwater sampling event for the 2<sup>nd</sup> Quarter 2009, the deviations from standard procedures of the GWP included the use of dedicated pumps in each of the wells and elimination of equipment rinsate and field blanks.





## 5.0 GROUNDWATER MONITORING RESULTS

Soil boring/well construction logs, historic data results, well numbering, and construction details are included in the previous work plans and quarterly sampling reports (see **Section 7.0**). Water level depths and groundwater field parameters for pH, temperature, conductivity, ORP, dissolved oxygen, and turbidity for the 2<sup>nd</sup> Quarter/June 2009 sampling event are shown on **Table 5-1**:

**Table 5-1- Summary of Field Parameters**

Sample ID	Date	Depth to Water <i>ft below TOC*</i>	Water Elevation <i>Feet msl</i>	Temperature <i>°C</i>	Specific Conductivity <i>uS/cm</i>
23L4MW01AW	6/29/2009	16.74	514.66	10.18	29
23L4MW01BW	6/29/2009	13.52	516.05	10.34	25
23L4MW02AW	6/29/2009	35.5	484.43	11.17	33
23L4MW02BW	6/29/2009	53.62	464.84	10.85	73
23L4MW03AW	6/26/2009	30.17	484.68	11.7	23
23L4MW03BW	6/26/2009	27.51	483.96	11.32	48
23L4MW04AW	6/29/2009	28.23	483.56	10.19	17
23L4MW05AW	6/26/2009	24.58	485.33	10.64	27
23L4MW07BW	6/26/2009	40.2	440.6	9.94	33
23L4MW17W	6/24/2009	10.88	350.6	11.89	220
23L4MW18W	6/24/2009	11.77	351.07	11.48	121
23LCMW01SW	6/26/2009	5.72	284.435	11.13	85
23LCMW01DW	6/26/2009	6.01	284.242	12.02	88
23LCMW02SW	6/26/2009	6.51	284.676	11.25	86
23LCMW02DW	6/26/2009	6.86	284.731	11.76	91
23LCMW03SW	6/26/2009	5.76	285.15	10.98	85
23LCMW03DW	6/26/2009	5.82	285.16	12.02	93
23LCMW04SW	6/24/2009	5.56	286.067	12.92	9.2
23LCMW04DW	6/24/2009	6.21	285.579	10.89	100

Sample ID	Date	Dissolved Oxygen <i>mg/l</i>	pH <i>S.U.</i>	Oxydation Reduction Potential <i>Millivolts</i>	Turbidity <i>NTU</i>
23L4MW01AW	6/29/2009	5.2	5.09	269.1	0.41
23L4MW01BW	6/29/2009	3.9	5.21	279.3	0.44
23L4MW02AW	6/29/2009	3.1	4.8	337.9	0.22
23L4MW02BW	6/29/2009	0.9	5.69	58.6	0.22
23L4MW03AW	6/26/2009	7.3	5.67	252.1	0.39
23L4MW03BW	6/26/2009	7.49	5.93	237	3.43
23L4MW04AW	6/29/2009	2.8	4.95	297.4	1.21
23L4MW05AW	6/26/2009	7.16	5.57	247.2	0.68
23L4MW07BW	6/26/2009	7.28	5.59	237	0.11
23L4MW17W	6/24/2009	5.62	7.52	84.7	4.01
23L4MW18W	6/24/2009	7.02	6.32	140.2	2.72
23LCMW01SW	6/26/2009	7.78	6.79	158.9	0.28
23LCMW01DW	6/26/2009	7.81	6.87	149.6	0.36
23LCMW02SW	6/26/2009	7.98	7.05	141.5	0.4
23LCMW02DW	6/26/2009	7.92	7	134.6	0.55
23LCMW03SW	6/26/2009	8.21	6.81	121.8	0.42
23LCMW03DW	6/26/2009	8.16	6.87	153.7	0.35
23LCMW04SW	6/24/2009	8.52	6.7	174.5	0.51
23LCMW04DW	6/24/2009	8.47	6.66	90.5	0.64

ft below TOC – Feet below top of casing  
 ft msl – feet to mean sea level  
 °C – degrees Celsius  
 uS/cm – microSiemens per centimeter

mg/L – milligram per liter/part per million  
 s.u. – standard unit  
 ntu - nephelometric turbidity units



## 5.1 Sentinel Wells

Groundwater samples were collected on June 24<sup>th</sup> and 25<sup>th</sup>, 2009 from the Sentinel Wells, which consist of four monitoring well pairs located at along the southwestern corner of CBMT where Lacamas Creek exits the site (see **Figures 1.2 and 3.5**). A field duplicate sample (labeled LCMW460W) was collected from Monitoring Well LCMW04S. Extra volume of groundwater was collected from Monitoring Well LCMW03B for the purpose of supplying extra water for laboratory MS/MSD samples. Trip blanks accompanied all groundwater VOC sample containers.

VOCs, explosive compounds, RDX, HMX, nitroglycerine, PETN, and picric acid were not detected in any of the Sentinel Wells groundwater samples; see **Table 5-2** below:

**Table 5-2  
 Summary of Sentinel Well Groundwater Sample Results**

Site Sample ID Sample Date/Time Units	23LCMW01DW 6/25/2009 ug/L	23LCMW01SW 6/25/2009 ug/L	23LCMW02DW 6/25/2009 ug/L	23LCMW02SW 6/25/2009 ug/L
1,1,1-Trichloroethane	ND	ND	ND	ND
1,1-Dichloroethane	ND	ND	ND	ND
1,1-Dichloroethene	ND	ND	ND	ND
1,3,5-Trinitrobenzene	ND	ND	ND	ND
2,4-Dinitrotoluene	ND	ND	ND	ND
2-Nitrotoluene	ND	ND	ND	ND
Dichlorodifluoromethane	ND	ND	ND	ND
HMX	ND	ND	ND	ND
Perchlorate	ND	ND	ND	ND
RDX	ND	ND	ND	ND
Toluene	ND	ND	ND	ND
Trichloroethene	ND	ND	ND	ND

Site Sample ID Sample Date/Time Units	23LCMW03DW 6/24/2009 ug/L	23LCMW03SW 6/25/2009 ug/L	23LCMW04DW 6/24/209 ug/L	23LCMW04SW 6/24/2009 ug/L
1,1,1-Trichloroethane	ND	ND	ND	ND
1,1-Dichloroethane	ND	ND	ND	ND
1,1-Dichloroethene	ND	ND	ND	ND
1,3,5-Trinitrobenzene	ND	ND	ND	ND
2,4-Dinitrotoluene	ND	ND	ND	ND
2-Nitrotoluene	ND	ND	ND	ND
Dichlorodifluoromethane	ND	ND	ND	ND
HMX	ND	ND	ND	ND
Perchlorate	ND	ND	ND	ND
RDX	ND	ND	ND	ND
Toluene	ND	ND	ND	ND
Trichloroethene	ND	ND	ND	ND

ug/L – microgram per liter/part per billion

ND – Non-detectable

## 5.2 Landfill 4/Demolition Area 1

Quarterly groundwater samples have been collected from eleven wells associated with LF4/DA1 (see **Figures 1.2 and 3.8**) since 2001, except between April 2002 and December 2003. Eight of those wells surround the former landfill area (LF4-MW01A&B, LF4-MW02A&B, LF4-MW03A&B, LF4-MW04A, and LF4-MW05A), with one located down/side gradient to the south (LF4-MW07B) and two wells (LF4-MW17 and LF4-MW18) located where the North Fork of Lacamas Creek enters the alluvial deposits in the valley (part of the Troutdale Sole Source Aquifer).

Naturally occurring levels of inorganic elements (metals) were observed in the parts per billion ranges in all of the wells (Ecology 1994). Historically the highest observed concentrations were generally associated with samples that had a high level of suspended solids in them. The comparable filtered sample for many of the metals have displayed lower concentrations, indicating the elevated reading was the result of the acid preservative dissolving materials from the soil particles suspended in the sample. In June 2005 the total mercury concentration in well L4MW03A (5.6 ug/L) exceeded the MTCA Method A criteria (2 ug/L). The associated dissolved mercury analysis (0.09 ug/L) was well below MTCA Method A criteria. All previous and subsequent mercury analyses from this well have been at non-detect levels (that are below the MTCA criteria). A significant fraction of the other site-wide wells also had laboratory reported mercury detections (all below MTCA criteria) in this single sampling event (LFMW1B, 2A, 2B, 4B, 4A, 7B, 17 18; MW3S, 3D, 4S, 4D). This pattern is interpreted to be a laboratory error or laboratory equipment contamination.

No other metals (total or dissolved) from RAU 2C have exceeded the MTCA Method A or Method B criteria. Ecology approved the removal of metals from the routine monitoring parameters at LF4/DA1 in 2006.

Groundwater samples were collected from the LF4/DA1 monitoring wells (see **Figures 1.2 and 3.8**) on June 25<sup>th</sup>, 26<sup>th</sup>, and 29<sup>th</sup>, 2009. A field duplicate sample (labeled L4MW460W) was collected from Monitoring Well L4MW03B. Trip blanks accompanied all groundwater VOC sample containers.

PETN, picric acid, and nitroglycerin, were not detected in any of the groundwater samples from shallow or deep monitoring wells. HMX and RDX were not detected in Monitoring Wells L4MW17, L4MW18, L4MW01B, and L4MW07B. HMX and RDX were both detected in monitoring wells L4MW02A, L4MW02B, L4MW03A, and L4MW05A. RDX only was detected in monitoring wells L4MW01A, L4MW03B, and L4MW04A.

Perchlorate was detected in groundwater samples from all the LF4/DA1 wells except L4MW17 and L4MW18. The concentrations range from 1.1 µg/L at L4MW01B to 431 µg/L at L4MW02B. The highest levels of HMX, RDX, and perchlorate were found in the groundwater samples from the paired monitoring wells MW02A and MW02B.

VOCs detected at well L4MW02B included 1,1-dichloroethane, 1,1-dichloroethene, 1,1,1-trichloroethane, 2,4-dinitrotoluene, and dichlorodifluoromethane, although below applicable MTCA Method B cleanup values. VOCs were not detected in any of the other monitoring wells; see **Table 5-3** below:

Complete copies of the laboratory data packages and validation reports are included electronically on the attached CD as **Appendix A**.



**Table 5-3**  
**Summary of LF4/DA1 Well Groundwater Sample Results**

Site Sample ID Sample Date/Time Units	23L4MW01AW 6/29/2009 ug/L	23L4MW01BW 6/29/2009 ug/L	23L4MW02AW 6/29/2009 ug/L	23L4MW02BW 6/29/2009 ug/L
1,1,1-Trichloroethane	ND	ND	ND	19
1,1-Dichloroethane	ND	ND	ND	19.2
1,1-Dichloroethene	ND	ND	ND	6.15
1,3,5-Trinitrobenzene	ND	ND	ND	ND
2,4-Dinitrotoluene	ND	ND	ND	0.36 J
2-Nitrotoluene	ND	ND	ND	ND
Dichlorodifluoromethane	ND	ND	ND	23.9
HMX	ND	ND	3.8 J	4.3 J
Perchlorate	2.9	1.1	195	431
RDX	0.13	ND	19 J	84 J
Toluene	ND	ND	ND	ND
Trichloroethene	ND	ND	ND	ND

Site Sample ID Sample Date/Time Units	23L4MW03AW 6/26/2009 ug/L	23L4MW03BW 6/26/2009 ug/L	23L4MW04AW 6/29/2009 ug/L	23LF4MW5A 6/26/2009 ug/L
1,1,1-Trichloroethane	ND	ND	ND	ND
1,1-Dichloroethane	ND	ND	ND	ND
1,1-Dichloroethene	ND	ND	ND	ND
1,3,5-Trinitrobenzene	ND	ND	ND	ND
2,4-Dinitrotoluene	ND	ND	ND	ND
2-Nitrotoluene	ND	ND	ND	ND
Dichlorodifluoromethane	ND	ND	ND	ND
HMX	0.38	ND	ND	0.25 J
Perchlorate	82.6	41.9	34.2	35.9
RDX	9.4	4.1	2.8 J	4.1 J
Toluene	ND	ND	ND	ND
Trichloroethene	ND	ND	ND	ND

Site Sample ID Sample Date/Time Units	23LF4MW7B 6/26/2009 ug/L	23LF4MW17SW 6/25/2009 ug/L	23LF4MW18SW 6/25/2009 ug/L
1,1,1-Trichloroethane	ND	ND	ND
1,1-Dichloroethane	ND	ND	ND
1,1-Dichloroethene	ND	ND	ND
1,3,5-Trinitrobenzene	ND	ND	ND
2,4-Dinitrotoluene	ND	ND	ND
2-Nitrotoluene	ND	ND	ND
Dichlorodifluoromethane	ND	ND	ND
HMX	ND	ND	ND
Perchlorate	2.3	ND	ND
RDX	ND	ND	ND
Toluene	ND	ND	ND
Trichloroethene	ND	ND	ND

ug/L – microgram per liter/part per billion

J – estimated value

ND – Non-detectable



## 6.0 GROUNDWATER QUALITY TRENDS

### 6.1 Sentinel Wells

The Sentinel Well groundwater result trends are as follows:

- With the use of dedicated pumps and low flow purging/sampling techniques (which obtain water samples with lower turbidity), the total and dissolved metals concentrations have decreased significantly. All of the total and dissolved metals detections in groundwater from these wells were below MTCA Method A and B regulatory screening levels.
- Petroleum hydrocarbons have not been detected in any of the Sentinel Wells throughout the monitoring, except for an isolated detection of diesel range petroleum hydrocarbons in LCMW02DW at 0.15 mg/L in January 2006.
- Perchlorate and explosive constituents have not been detected above the laboratory detection limits at the Sentinel Wells, following data validation and data quality assessments.

### 6.2 LF4/DA1Wells – Data Trends

Based on our recent review of historic groundwater data, the following observations summarize groundwater conditions at and around the LF4/DA1 site; detections are shown graphically for each well in **Appendix B**.

- All of the VOCs detected (primarily at LF4-MW-2B) continue to be well below MTCA Method A and B Cleanup Levels (see **Table 6-1**). Concentrations of 1,1-dichloroethene, 1,1-dichloroethane and 1,1,1-trichloroethane have been decreasing slowly and dichlorodifluoromethane, 1,2-dichloroethane, tetrachloroethene and trichloroethene results have been non-detectable for the last year.
- HMX and RDX concentrations in groundwater have been either stable, below MTCA levels, or decreasing slowly with consistent concentration distributions throughout all of the 27 LF4/DA1 groundwater sampling events (2001 to 2009; see **Figure 6.1**).
- Well LF4-MW-1A – the shallow upgradient well perchlorate concentrations have decreased to previous levels (ranging from 1.6 to 7 µg/L) from the 36 µg/L detected in 4th Quarter 2008 during and following a heavy precipitation event and correlate with RDX variations (see **Figure 6.2**).
- Well LF4-MW-1B – the deep upgradient well has low perchlorate concentrations; the adjacent shallow well LF-MW-1A has a history of low perchlorate and RDX concentrations. Since this well is located upgradient of the LF4/DA1, neither the detection nor absence of perchlorate at this well affects the monitoring program.

- Well LF-MW-2A – perchlorate concentrations appear to have reached a degree of equilibrium since 2005 with a consistent concentrations of perchlorate, RDX and HMX, which are all clearly affected by seasonal changes in recharge (see **Figure 6.3**). The seasonal variation appears to be inversely correlated with increased precipitation/groundwater elevations:
  - The lowest reported perchlorate concentrations, the highest measured groundwater levels and the lowest reported precipitation totals have generally occurred in the first quarter (Winter) events each year.
  - The highest reported perchlorate concentrations, the lowest measured groundwater levels, and the highest reported precipitation totals usually occurred in the third quarter (Summer) events of each year.
- L4MW2B perchlorate levels follow a stable, quasi-seasonal pattern. The 1<sup>st</sup> and 2<sup>nd</sup> Quarters in 2009 show the beginning of the downward portion of this pattern. The clearly seasonal perchlorate/RDX/HMX concentration patterns observed in L4-MW-2A are not repeated in the LF-MW-2B data (see **Figure 6.4**).
- Well L4-MW-3A - perchlorate concentrations have remained relatively stable with a slightly decreasing trend since a peak concentration of 120 µg/L occurred during the 3<sup>rd</sup> quarter 2006 sampling event. RDX concentrations are stable at about 10 ug/L.
- Well L4-MW-3B - perchlorate concentrations have remained relatively stable with an overall decreasing trend since a peak concentration (55 µg/L) was observed in the 3<sup>rd</sup> and 4<sup>th</sup> quarter 2006 sampling events. RDX concentrations are stable at 5 ug/L.
- Well L4-MW-4A - perchlorate concentrations have remained relatively stable (29 to 34 µg/L) since a peak concentration (40 µg/L) was observed in the 4<sup>th</sup> quarter 2006 and 2<sup>nd</sup> quarter 2007 sampling events. The common laboratory contaminant, methylene chloride that was detected (0.14 ppb) in the 2<sup>nd</sup> Quarter 2008 sample, has not been detected in later events.
- Well L4-MW-5A - perchlorate concentrations have been generally decreasing from a peak of 64 ppb in the initial sampling event in the 3<sup>rd</sup> quarter 2001 to less than 40 ppb during the last 7 quarters. The trace detections of tetrachloroethene have been non-detectable for the last 3 quarters. RDX concentrations have been consistently less than 5 ug/L.
- Well L4-MW-7B - perchlorate concentrations have been generally stable at 2 to 3 ppb for the last 20 quarterly sampling events; with the exception of an apparent field cross contamination issue during the 1<sup>st</sup> quarter 2006 event.
- Well L4-MW-17 – the 2<sup>nd</sup> Quarter 2008 estimated (above the MDL but below the MRL) concentrations of 1,2,4-Trimethylbenzene and Naphthalene (0.12 and 0.35 µg/L, respectively) have not been detected in subsequent events.



- Well L4-MW-17 and 18 – the monitoring wells located at the beginning of the Central Valley Floor were non-detectable for perchlorate throughout the 20 sampling events

**Table 6-1**  
**Summary of Maximum Groundwater Monitoring Detections vs MCTA Cleanup Levels**

Contaminant	Concentration (ppb/μg/L)		MCTA Cleanup Level Groundwater (ppb/μg/L)	
	Maximum Detection <sup>(1)</sup>	Well ID	Method A	Method B
HMX	4.6	L4MW02A		400
<b>RDX</b>	<b>120</b>	<b>L4MW02B</b>		<b>0.8</b>
<b>Perchlorate</b>	<b>530</b>			<b>11</b>
Picric Acid	2.9	L4MW02B	(2)	
Nitroglycerine	8.4		(2)	
1,1-Dichloroethene	36			400
1,1- Dichloroethane	45			1600
1,1,1-Trichloroethane	>200 in '03 >50 in '07			200
1,2-Dichloroethane	1.58			5 <sup>(2)</sup>
2,4-Dinitrotoluene	0.36	L4MW02B	(2)	32
Isopropylbenzene	0.2		(2)	
Methylene Chloride	2.58		5 <sup>(2)</sup>	
n-Propylbenzene	0.2		(2)	48
Tetrachloroethene	1.1		5	
Trichloroethene	0.26		5 <sup>(2)</sup>	
Trichlorofluoromethane	0.22		5 <sup>(2)</sup>	2,400
Dichlorodifluoromethane	190			1,600
1,3,5 Trinitrobenzene	0.13	L4MW01A		43,000
Naphthalene	0.35	L4MW17	160 <sup>(2)</sup>	

<sup>(1)</sup> **Bold** values exceeded MCTA A or B Cleanup Levels.

<sup>(2)</sup> Not retained due to low frequency detection (one or two times out of > 250 samples).

The quarterly sample results continue to be less than the Applicable Cleanup Levels as presented in the Draft RI/FS for Site-Wide Groundwater (RAU2C/BCRRT August 2009) currently under review by Ecology.

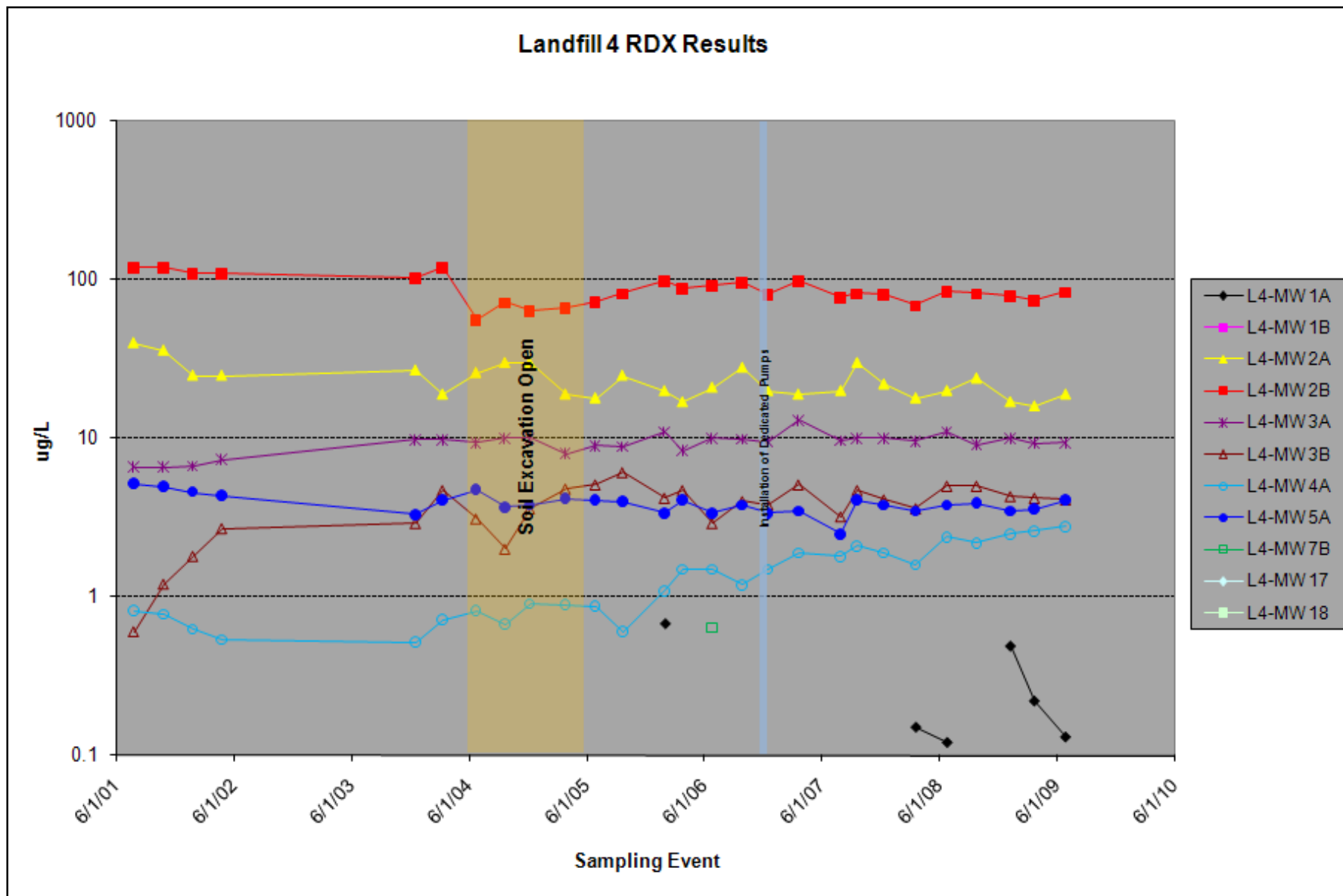


Figure 6.1 LF4/DA1 Groundwater Monitoring RDX Results



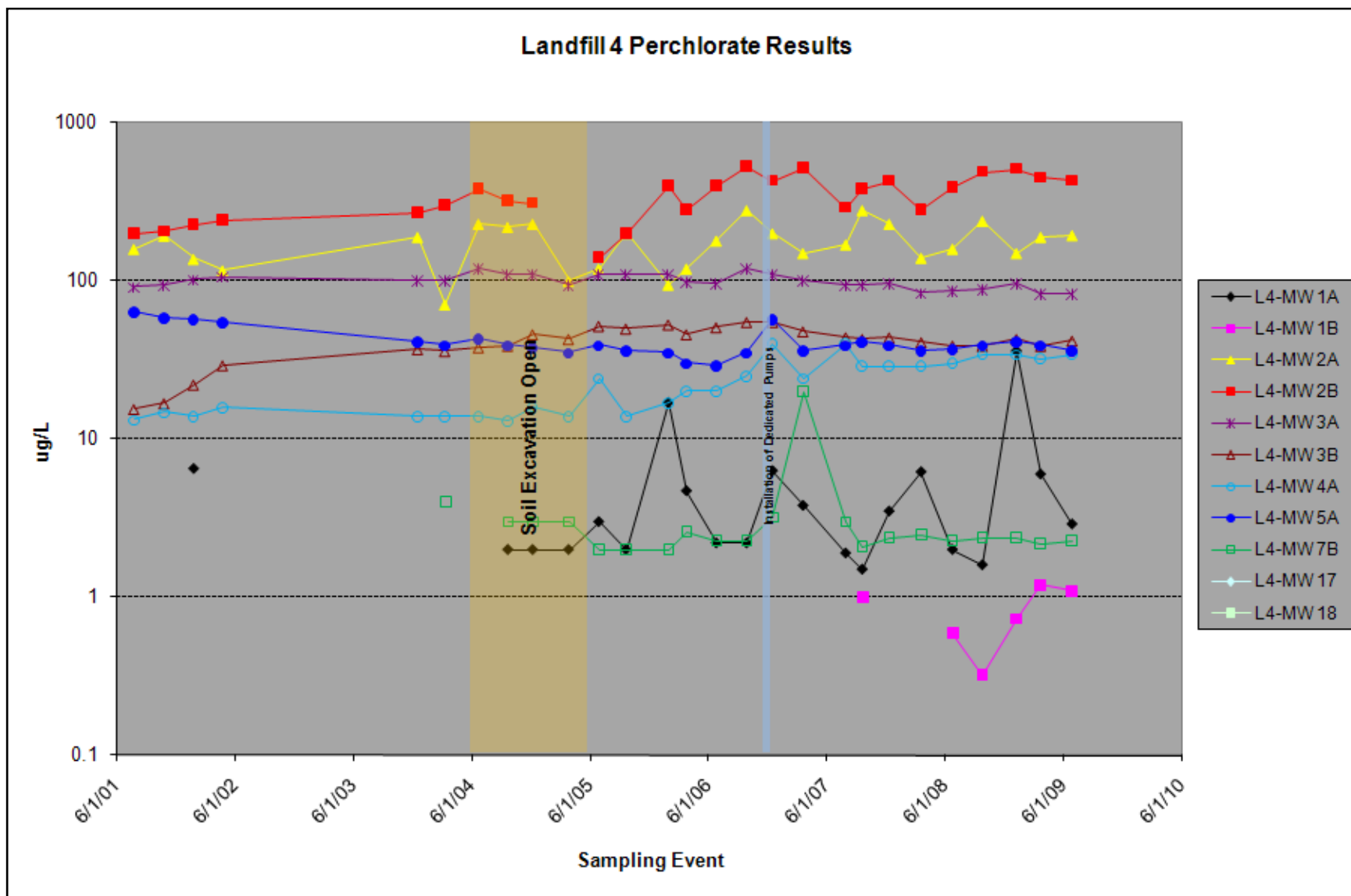
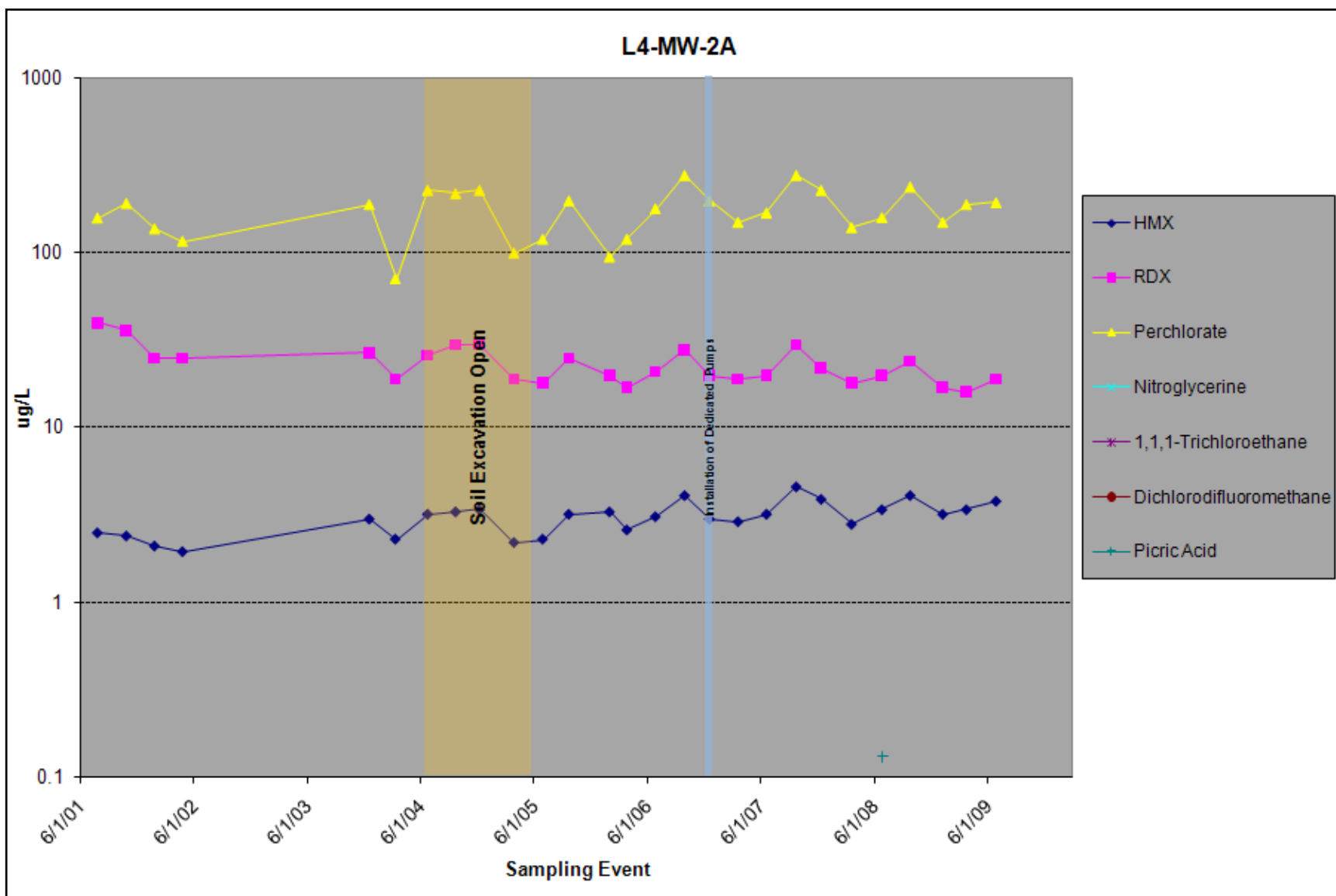


Figure 6.2 LF4/DA1 Groundwater Monitoring Perchlorate Results



**Figure 7.3** LF4/DA1 Perchlorate Results, Groundwater Elevation, and Precipitation - LF4-MW-2A

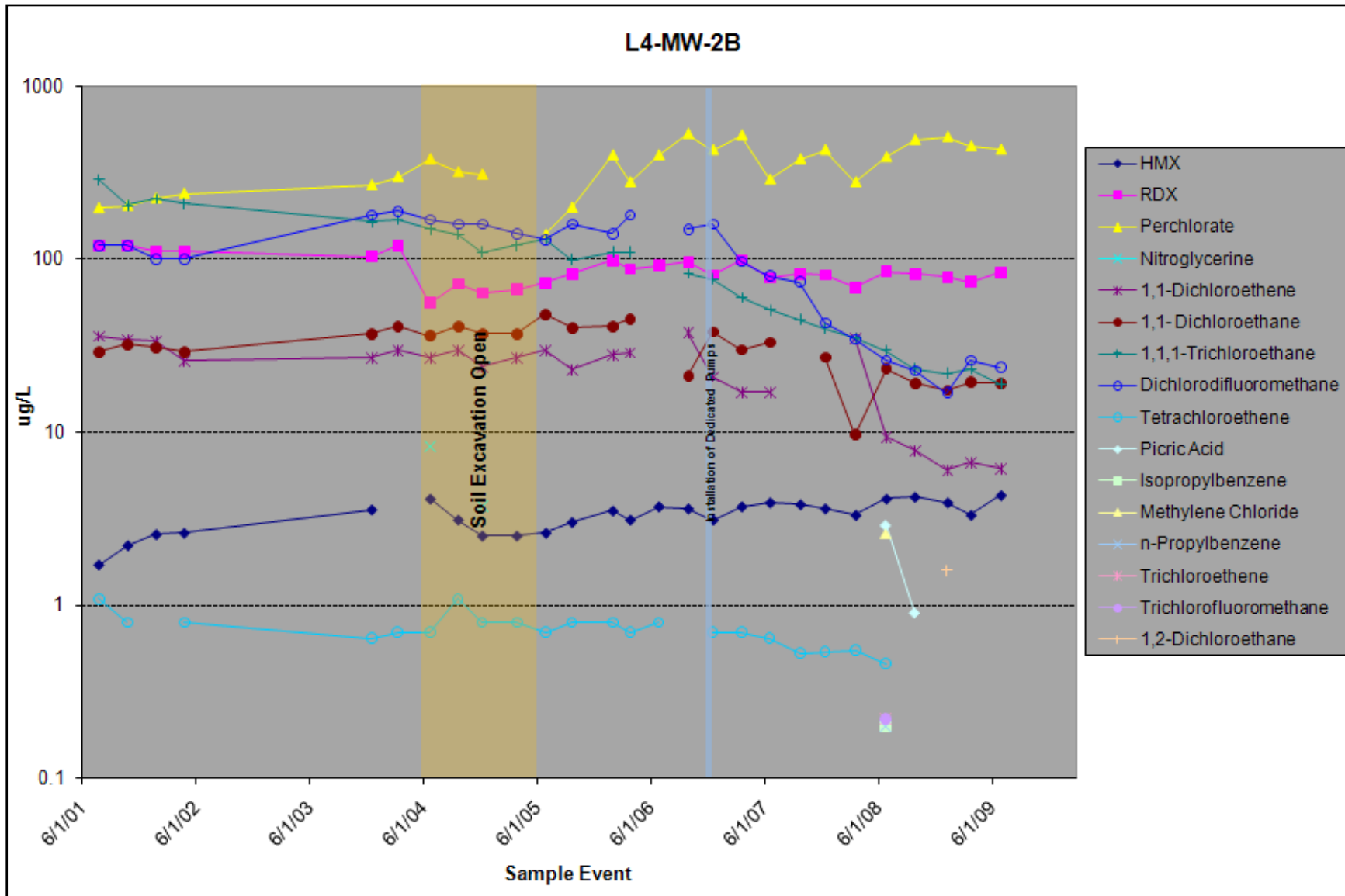


Figure 6.4 LF4/DA1 Perchlorate Results, Groundwater Elevation, and Precipitation - LF4-MW-2B



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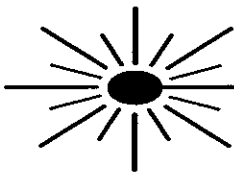
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**Data Validation Report**

**Michael Baker Jr., Inc**

**Camp Bonneville**

**SDG#: PSF0784**

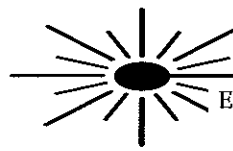


**ECT.CON INC.**

Environmental and Computer  
Technology Consultants

3531 Fox Chase Drive  
Imperial, PA 15126  
(724) 695-8042  
FAX (724) 695-2698  
e-mail: [ectconinc@comcast.net](mailto:ectconinc@comcast.net)





## Data Validation Report

SDG#	PSF0784
Validation Report Date	August 17, 2009
Validation Guidance	USEPA CLP National Functional Guidelines for Data Review
Client Name	Michael Baker Jr., Inc.
Project Name	Camp Bonneville
Laboratory	TestAmerica
Analytical Parameters	Explosives
Analytical Method	SW-846 8330

### Samples/Matrix:

Date Sampled	Location	Portland ID	Sacramento ID	Explosives	Matrix
06/24/09	23LCMW04DW	PSF0784-01	LFQAK	X	Aqueous
06/24/09	23LCMW04SW	PSF0784-02	LFQAL	X	Aqueous
06/24/09	23LCMW0460W	PSF0784-03	LFQAM	X	Aqueous

Analytical data in this report were screened to determine analytical limitations of the data based on specific quality control criteria. This screening assumes analytical results are correct as reported and merely provides an interpretation of the reported quality control results. Laboratory calculations have been verified as part of this validation. Specific findings on analytical limitations are presented in this report. Annotated Form 1s or spreadsheets for samples reviewed are included after the Data Assessment Findings. Form 1s for the MS/MSD samples and spreadsheets are not annotated.

### SUMMARY

The sample set for the Camp Bonneville site consists of three aqueous field samples. These samples were analyzed for explosives as provided in the Table 1. The findings presented in this review of the analytical data assume that the information presented by the analytical laboratory is correct.

The explosives findings are based upon the assessment of the following:

- \* ● Data Completeness
  - \* ● Holding Times
  - \* ● Calibration (Initial and Continuing)
  - Blanks
  - \* ● System Monitoring Compounds (Surrogate Spikes)
  - \* ● Matrix Spike/Matrix Spike Duplicates
  - \* ● Laboratory Control Standard Results
  - \* ● Target Compound Identification
  - \* ● Compound Quantification and Reported Contract Quantitation Limits
  - \* ● System Performance
- \* Criteria were met for this evaluation item.

This evaluation was conducted in accordance with USEPA CLP National Functional Guidelines for Organic Data Review and the analytical method. Findings from this evaluation should be considered when using the analytical data. This report presents a summary of the data qualifications based on the review of the aforementioned evaluation criteria. This is followed by annotated Form 1s/ spreadsheets. Finally, the worksheets used to perform the evaluation are provided.

## **FINDINGS**

### **1. Blanks**

The laboratory method blank exhibited contamination for the following parameter:

Blank	Compound	Maximum Concentration (ppb)	Action Limit (ppm)	Action
LFW791	1,3,5-Trinitrobenzene	0.038	0.19	U sample results < RL

RL - reporting limit

## **NOTES**

### **Matrix Spike/Matrix Spike Duplicate Results**

A MS/MSD was not analyzed with this SDG. A laboratory control sample was analyzed. Data were not qualified on this basis.

### **Laboratory Control Sample Results**

Recovery of 12 out of 17 compounds exceeded the laboratory's upper quality control limits. The affected compounds were not detected in the associated samples. Data were not qualified on this basis.

### **Field Duplicate Results**

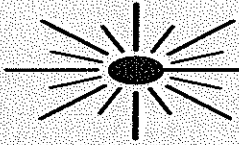
No field duplicates were included with this SDG. Data were not qualified on this basis.

  
Data Reviewer

8/12/09  
Date

### **Glossary of Data Qualifiers**

U	Not Detected.	The associated number indicates approximate sample concentration necessary to be detected.
UJ	Not Detected.	Quantitation limit may be inaccurate or imprecise.
J	Analyte Present.	Reported value may not be accurate or precise.
N	Consider Present.	Tentative identification. Special methods may be needed to confirm its presence or absence in future sampling efforts.
R	Unusable Result.	Analyte may or may not be present in the sample.
UR	Unusable Result.	Analyte may or may not be present in the sample.



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*Annotated Form 1's*  
*(Spreadsheet)*

TestAmerica Portland

Client Sample ID: PSF0784-01

HPLC *23LCM.W040W*

Lot-Sample #...: G9F260326-001 Work Order #...: LFQAK1AC Matrix.....: WATER  
 Date Sampled...: 06/24/09 Date Received...: 06/26/09  
 Prep Date.....: 07/01/09 Analysis Date...: 07/02/09  
 Prep Batch #...: 9182192  
 Dilution Factor: 1 Method.....: SW846 8330

PARAMETER	RESULT	REPORTING		
		LIMIT	UNITS	MDL
Nitroglycerin	ND	0.65	ug/L	0.15
PETN	ND	0.65	ug/L	0.23
Picric Acid	ND	1.0	ug/L	0.12
2-Amino-4,6-dinitrotoluene	ND	0.20	ug/L	0.10
4-Amino-2,6-dinitrotoluene	ND	0.10	ug/L	0.022
1,3-Dinitrobenzene	ND	0.10	ug/L	0.050
2,4-Dinitrotoluene	ND	0.10	ug/L	0.050
2,6-Dinitrotoluene	ND	0.10	ug/L	0.050
HMX	ND	0.10	ug/L	0.027
Nitrobenzene	ND	0.10	ug/L	0.050
2-Nitrotoluene	ND	0.40	ug/L	0.072
3-Nitrotoluene	ND	0.40	ug/L	0.062
4-Nitrotoluene	ND	0.50	ug/L	0.072
RDX	ND	0.10	ug/L	0.065
Tetryl	ND	0.10	ug/L	0.050
1,3,5-Trinitrobenzene	<i>0.036 J, B</i>	0.10	ug/L	0.031
2,4,6-Trinitrotoluene	ND	0.10	ug/L	0.024

SURROGATE	PERCENT	RECOVERY
	RECOVERY	LIMITS
3,4-Dinitrotoluene	104	(79 - 111)

NOTE(S):

- J Estimated result. Result is less than RL.
- B Method blank contamination. The associated method blank contains the target analyte at a reportable level.

*UK 8/6/09*



TestAmerica Portland

Client Sample ID: PSF0784-02

HPLC

23LCM045W

Lot-Sample #...: G9F260326-002 Work Order #...: LFQAL1AC Matrix.....: WATER  
 Date Sampled...: 06/24/09 Date Received...: 06/26/09  
 Prep Date.....: 07/01/09 Analysis Date...: 07/02/09  
 Prep Batch #...: 9182192  
 Dilution Factor: 1 Method.....: SW846 8330

PARAMETER	RESULT	REPORTING		
		LIMIT	UNITS	MDL
Nitroglycerin	ND	0.65	ug/L	0.15
PETN	ND	0.65	ug/L	0.23
Picric Acid	ND	1.0	ug/L	0.12
2-Amino-4,6-dinitrotoluene	ND	0.20	ug/L	0.10
4-Amino-2,6-dinitrotoluene	ND	0.10	ug/L	0.022
1,3-Dinitrobenzene	ND	0.10	ug/L	0.050
2,4-Dinitrotoluene	ND	0.10	ug/L	0.050
2,6-Dinitrotoluene	ND	0.10	ug/L	0.050
HMX	ND	0.10	ug/L	0.027
Nitrobenzene	ND	0.10	ug/L	0.050
2-Nitrotoluene	ND	0.40	ug/L	0.072
3-Nitrotoluene	ND	0.40	ug/L	0.062
4-Nitrotoluene	ND	0.50	ug/L	0.072
RDX	ND	0.10	ug/L	0.065
Tetryl	ND	0.10	ug/L	0.050
1,3,5-Trinitrobenzene	0.033 J,B	0.10	ug/L	0.031
2,4,6-Trinitrotoluene	ND	0.10	ug/L	0.024

SURROGATE	PERCENT	RECOVERY
	RECOVERY	LIMITS
3,4-Dinitrotoluene	105	(79 - 111)

NOTE(S):

- J Estimated result. Result is less than RL.
- B Method blank contamination. The associated method blank contains the target analyte at a reportable level

WJ  
8/2/09

TestAmerica Portland

Client Sample ID: PSF0784-03

HPLC

23 LEM W0460W

Lot-Sample #...: G9F260326-003    Work Order #...: LFQAM1AC    Matrix.....: WATER  
 Date Sampled...: 06/24/09    Date Received...: 06/26/09  
 Prep Date.....: 07/01/09    Analysis Date...: 07/02/09  
 Prep Batch #...: 9182192  
 Dilution Factor: 0.99    Method.....: SW846 8330

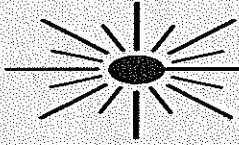
PARAMETER	RESULT	REPORTING LIMIT	UNITS	MDL
Nitroglycerin	ND	0.64	ug/L	0.15
PETN	ND	0.64	ug/L	0.23
Picric Acid	ND	0.99	ug/L	0.12
2-Amino-4,6-dinitrotoluene	ND	0.20	ug/L	0.099
4-Amino-2,6-dinitrotoluene	ND	0.099	ug/L	0.022
1,3-Dinitrobenzene	ND	0.099	ug/L	0.050
2,4-Dinitrotoluene	ND	0.099	ug/L	0.050
2,6-Dinitrotoluene	ND	0.099	ug/L	0.050
HMX	ND	0.099	ug/L	0.027
Nitrobenzene	ND	0.099	ug/L	0.050
2-Nitrotoluene	ND	0.40	ug/L	0.071
3-Nitrotoluene	ND	0.40	ug/L	0.061
4-Nitrotoluene	ND	0.50	ug/L	0.071
RDX	ND	0.099	ug/L	0.064
Tetryl	ND	0.099	ug/L	0.050
1,3,5-Trinitrobenzene	<del>0.035</del> J,B	0.099	ug/L	0.031
2,4,6-Trinitrotoluene	ND	0.099	ug/L	0.024
	PERCENT	RECOVERY		
SURROGATE	RECOVERY	LIMITS		
3,4-Dinitrotoluene	106	(79 - 111)		

NOTE(S):

J Estimated result. Result is less than RL.

B Method blank contamination. The associated method blank contains the target analyte at a reportable level.

UK  
8/6/09



ECT.CON INC.

Environmental and Computer  
Technology Consultants

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## *Support Documentation*

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## *Laboratory Case Narrative*

## DATA DELIVERABLES PACKAGE

Michael Baker Jr., Inc.  
James D. Peyton  
5261 Fountain Drive, Suite A  
Crown Point, IN 46307

Client Project: Camp Bonneville Groundwater  
Client Project Number: 110002 Task 6200

TA Work Order#: PSF0784  
TA Project Manager: Estella Rieben

The total number of pages contained in this data package is:

139

July 31, 2009

TestAmerica - Portland  
9405 S. W. Nimbus Avenue  
Beaverton, Oregon 97008  
(503) 906-9200  
(503) 906-9210

→ 200s between columns  
for sample results  
→ field dupes





**Michael Baker Jr., Inc.**  
5261 Fountain Drive, Suite A  
Crown Point, IN 46307

Project Name: **Camp Bonneville Groundwater**  
Project Number: 110002 Task 6200  
Project Manager: James D. Peyton

Report Created:  
07/21/09 08:41

## ANALYTICAL REPORT FOR SAMPLES

Sample ID	Laboratory ID	Matrix	Date Sampled	Date Received
23LCMW04DW	PSF0784-01	Water	06/24/09 13:00	06/24/09 17:30
23LCMW04SW	PSF0784-02	Water	06/24/09 15:45	06/24/09 17:30
23LCMW0460W	PSF0784-03	Water	06/24/09 11:00	06/24/09 17:30
TB261	PSF0784-04	Water	06/24/09 00:00	06/24/09 17:30

TestAmerica Portland



Estella Rieben, Project Manager

*The results in this report apply to the samples analyzed in accordance with the chain of custody document. This analytical report shall not be reproduced except in full, without the written approval of the laboratory.*



# TestAmerica

THE LEADER IN ENVIRONMENTAL TESTING

July 28, 2009

**TestAmerica Project Number: G9F260326**

PO/Contract: PSF0784

Estella Rieben  
TestAmerica Portland  
Nimbus Corporate Center  
9405 SW Nimbus Ave  
Beaverton, OR 97008

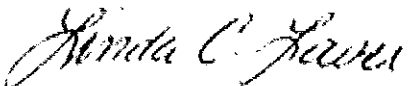
Dear Ms. Rieben,

This report contains the analytical results for the samples received under chain of custody by TestAmerica on June 26, 2009. These samples are associated with your Michael Baker Jr., Inc. project.

The test results in this report meet all NELAC requirements for parameters that accreditation is required or available. Any exceptions to NELAC requirements are noted in the case narrative. The case narrative is an integral part of this report.

If you have any questions, please feel free to call me at (916) 374-4362.

Sincerely,



Linda C. Laver  
Project Manager

## Case Narrative

### TestAmerica West Sacramento Project Number G9F260326

#### **WATER, 8330, Nitroaromatics & Nitramines**

Samples: 1, 2, 3

The laboratory control sample (LCS) associated with this extraction batch has percent recoveries for most analytes above the established control limits indicating a potential high bias in the data. These samples do not have detected concentrations above the reporting limit for these analytes and there is no adverse impact upon the data.

The matrix spike/matrix spike duplicate (MS/MSD) associated with this extraction batch also has recoveries and/or precision is outside the established control limits for many analytes. This anomaly is most likely matrix related.

There are no other anomalies associated with this project.

## Sample Summary

### TestAmerica West Sacramento Project Number G9F260326

<u>WO#</u>	<u>Sample #</u>	<u>Client Sample ID</u>	<u>Sampling Date</u>	<u>Received Date</u>
LFQAK	1	PSF0784-01	6/24/2009 01:00 PM	6/26/2009 09:00 AM
LFQAL	2	PSF0784-02	6/24/2009 03:45 PM	6/26/2009 09:00 AM
LFQAM	3	PSF0784-03	6/24/2009 11:00 AM	6/26/2009 09:00 AM

#### Notes(s):

- The analytical results of the samples listed above are presented on the following pages.
- All calculations are performed before rounding to avoid round-off errors in calculated results.
- Results noted as "ND" were not detected at or above the stated limit.
- This report must not be reproduced, except in full, without the written approval of the laboratory.
- Results for the following parameters are never reported on a dry weight basis: color, corrosivity, density, flashpoint, ignitability, layers, odor, paint filter test, pH, porosity, pressure, reactivity, redox potential, specific gravity, spot tests, solids, solubility, temperature, viscosity, and weight.



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## *Chain of Custody*

# TestAmerica

THE LEADER IN ENVIRONMENTAL TESTING

11720 North Creek Pkwy N Suite 401, Bothell, WA 98011 8244  
 11922 E. 1st Ave, Spokane, WA 99206 5292  
 5000 N. N. Highway, Vancouver, OR 97060 7143  
 100 W. Main Street, Portland, OR 97201 3333

253-270-9200 FAX 253-9210  
 509-325-9200 FAX 509-325-9200  
 503-946-9200 FAX 503-946-9200  
 503-253-9200 FAX 503-253-9200

## CHAIN OF CUSTODY REPORT

Work Order #: **P50784**

CLIENT: Michael Baker Corp		INVOICE TO: P&H				
REPORT TO: Margaret James		TURNAROUND REQUEST				
ADDRESS: 10000 N. 4th St, Suite A		in Business Days				
PHONE: 314-997-7363 FAX: 314-997-7363		Organic & Inorganic Analysis				
PROJECT NAME: Camp Brumville		<input type="checkbox"/> 10 <input type="checkbox"/> 11 <input type="checkbox"/> 12 <input type="checkbox"/> 13 <input type="checkbox"/> 14 <input type="checkbox"/> 15 <input type="checkbox"/> 16 <input type="checkbox"/> 17 <input type="checkbox"/> 18 <input type="checkbox"/> 19 <input type="checkbox"/> 20				
PROJECT NUMBER: 110000 T WACC		<input type="checkbox"/> 21 <input type="checkbox"/> 22 <input type="checkbox"/> 23 <input type="checkbox"/> 24 <input type="checkbox"/> 25 <input type="checkbox"/> 26 <input type="checkbox"/> 27 <input type="checkbox"/> 28 <input type="checkbox"/> 29 <input type="checkbox"/> 30				
SAMPLED BY: Margaret James		OTHER Specify				
CLIENT SAMPLE IDENTIFICATION		* Turnaround Requests less than standard may incur Rush Charges				
SAMPLING DATE/TIME		MATRIX # OF CONT. LOCATION/ COMMENTS TA W/O ID				
1. 230CMW04DW	6/24/09 - 13:00	W	6	NOTICS		
2. 230CMW04SW	6/24/09 - 1545	W	6			
3. 230CMW04DW	6/24/09 - 1100	W	6			
4. TD261	NA					
5.						
6.						
7.						
8.						
9.						
10.						
RELEASED BY: McDann	DATE: 6/24/09	RECEIVED BY: Bob F...	DATE: 6/24/09	FIRM: TAP	DATE: 6/24/09	TIME: 17:30
PRINT NAME: Margaret James	TIME: 16:00	PRINT NAME: Bob F...	TIME: 16:00	FIRM: TAP	DATE: 6/24/09	TIME: 17:30
RELEASED BY: Bob F...	DATE: 6/24/09	RECEIVED BY: Margaret James	DATE: 6/24/09	FIRM: TAP	DATE: 6/24/09	TIME: 17:30
PRINT NAME: Bob F...	TIME: 17:30	PRINT NAME: Margaret James	TIME: 17:30	FIRM: TAP	DATE: 6/24/09	TIME: 17:30
ADDITIONAL REMARKS:		TEMP: 9.0		PAGE 4.0		OF

TestAmerica Portland  
**Sample Receiving Checklist**

Work Order #: PSFO784 Date/Time Received: 6/24/09 1730  
 Client Name and Project: Michael Baker Camp Bonneville

Time Zone:  
 EDT/EST     CDT/CST     MDT/MST     PDT/PST     AK     OTHER

**Unpacking Checks:**

Cooler #(s): \_\_\_\_\_  
 Temperatures: \_\_\_\_\_  
 Digi #1    Digi #2    IR Gun  
                 (  Plastic  Glass )

**Temperature out of Range:**

\_\_\_\_ Not enough or No Ice  
 \_\_\_\_ Ice Melted  
 \_\_\_\_ W/in 4 Hrs of collection  
 \_\_\_\_ Other: \_\_\_\_\_

Initials: PS

- | N/A                                 | Yes                                 | No                       |  |
|-------------------------------------|-------------------------------------|--------------------------|--|
| <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/> | 1. If ESI client, were temp blanks received? If no, document on NOD.   |
| <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/> | 2. Cooler Seals intact? (N/A if hand delivered) if no, document on NOD.  |
|                                     | <input checked="" type="checkbox"/> | <input type="checkbox"/> | 3. Chain of Custody present? If no, document on NOD.   |
|                                     | <input checked="" type="checkbox"/> | <input type="checkbox"/> | 4. Bottles received intact? If no, document on NOD.  |
|                                     | <input checked="" type="checkbox"/> | <input type="checkbox"/> | 5. Sample is not multiphasic? If no, document on NOD.  |
|                                     | <input checked="" type="checkbox"/> | <input type="checkbox"/> | 6. Proper Container and preservatives used? If no, document on NOD.  |
| <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/> | 7. pH of all samples checked and meet requirements? If no, document on NOD.  |
| <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/> | 8. Cyanide samples checked for sulfides and meet requirements? If no, notify PM.   |
| <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/> | 9. HF Dilution required?   |
|                                     | <input checked="" type="checkbox"/> | <input type="checkbox"/> | 10. Sufficient volume provided for all analysis? If no, document on NOD and consult PM before proceeding.  |
|                                     | <input checked="" type="checkbox"/> | <input type="checkbox"/> | 11. Did chain of custody agree with samples received? If no, document on NOD.  |
|                                     | <input checked="" type="checkbox"/> | <input type="checkbox"/> | 12. Is the "Sampled by" section of the COC completed?  |
| <input type="checkbox"/>            | <input checked="" type="checkbox"/> | <input type="checkbox"/> | 13. Were VOA/Oil Syringe samples without headspace?  |
| <input type="checkbox"/>            | <input checked="" type="checkbox"/> | <input type="checkbox"/> | 14. Were VOA vials preserved? <input checked="" type="checkbox"/> HCl <input type="checkbox"/> Sodium Thiosulfate <input type="checkbox"/> Ascorbic Acid |
|                                     | <input checked="" type="checkbox"/> | <input type="checkbox"/> | 15. Did samples require preservation with sodium thiosulfate?  |
| <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/> | 16. If yes to #14, was the residual chlorine test negative? If no, document on NOD.  |
| <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/> | 17. Are dissolved/field filtered metals bottles sediment-free? If no, document on NOD.   |
| <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/> | 18. Is sufficient volume provided for client requested MS/MSD or matrix duplicates? If no, document on NOD and contact PM before proceeding.             |
| <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/> | 19. Are analyses with short holding times received in hold?  |
|                                     | <input checked="" type="checkbox"/> | <input type="checkbox"/> | 20. Was Standard Turn Around (TAT) requested?  |
|                                     | <input checked="" type="checkbox"/> | <input type="checkbox"/> | 21. Receipt date(s) < 48 hours past the collection date(s)? If no, notify PM.  |

TestAmerica Portland  
Sample Receiving Checklist

Work Order #: PSFO784

Login Checks:

Initials: PS

- | N/A                                 | Yes                                 | No                       |   |
|-------------------------------------|-------------------------------------|--------------------------|---|
|                                     | <input checked="" type="checkbox"/> | <input type="checkbox"/> | 22. Sufficient volume provided for all analysis? If no, document on NOD & contact PM.                                   |
| <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/> | 23. Sufficient volume provided for client requested MS/MSD or matrix duplicates? If no, document on NOD and contact PM. |
|                                     | <input checked="" type="checkbox"/> | <input type="checkbox"/> | 24. Did the chain of custody include "received by" and "relinquished by" signatures, dates and times?                   |
| <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/> | 25. Were special log in instructions read and followed?   |
|                                     | <input checked="" type="checkbox"/> | <input type="checkbox"/> | 26. Were tests logged checked against the COC?  |
| <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/> | 27. Were rush notices printed and delivered?  |
| <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/> | 28. Were short hold notices printed and delivered?  |
|                                     | <input checked="" type="checkbox"/> | <input type="checkbox"/> | 29. Were subcontract COCs printed?  |
| <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/> | 30. Was HF dilution logged?   |

Labeling and Storage Checks:

Initials: PS

- | N/A                                 | Yes                                 | No                       |   |
|-------------------------------------|-------------------------------------|--------------------------|---|
| <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/> | 31. Were the subcontracted samples/containers put in Sx fridge?                                     |
| <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/> | 32. Were sample bottles and COC double checked for dissolved/filtered metals?                       |
|                                     | <input checked="" type="checkbox"/> | <input type="checkbox"/> | 33. Did the sample ID, Date, and Time from label match what was logged?                             |
| <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/> | 34. Were Foreign sample stickers affixed to each container and containers stored in foreign fridge? |
| <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/> | 35. Were HF stickers affixed to each container, and containers stored in Sx fridge?                 |
| <input checked="" type="checkbox"/> | <input type="checkbox"/>            | <input type="checkbox"/> | 36. Was an NOD for created for noted discrepancies and placed in folder?                            |

Document any problems or discrepancies and the actions taken to resolve them on a Notice of Discrepancy form (NOD).





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# *Worksheets*

**HOLDING TIMES**

SAMPLE DATE	SAMPLE ID	Portland LAB ID	Sacramento Lab ID	Matrix	Prep	Anal
06/24/09	23LCMW04DW	PSF0784-01	LFQAK	Aq	07/01/09	7/2/09 IX
06/24/09	23LCMW04SW	PSF0784-02	LFQAL	Aq	07/01/09	7/2/09 IX
06/24/09	23LCMW0460W	PSF0784-03	LFQAM	Aq	07/01/09	7/2/09 IX

7 days from collection to extraction  
 Sample Date 6/24/2009 Extract By 7/1/2009  
 40 days from extraction to analysis  
 Sample Date 7/1/2009 Extract By 8/10/2009

**TARGET COMPOUNDS AND QUANTITATION LIMITS**

Was a target compound list provided by the client? No  
 Did Sample Form Is match the target compound list NA  
 Were required quantitation limits provided by the client? No  
 Did all compounds meet the required quantitation limits? NA

**SYSTEM MONITORING COMPOUNDS**

SAMPLE	SURROGATE	COLUMN I	DF	ACTION
ALL IN				

Were surrogate RTs within windows established by the ICAL? Yes  
 Were there any transcription errors between the raw data and Form 2? No  
 Were laboratory acceptance limits used as the basis for validation? 30-150  
 Did the laboratory provide CLP Form II or equivalent? Yes

04DW

SURR	AMOUNT FOUND	AMOUNT SPIKED	% R	FI
34-DNT A	2.607	2.5	104.3	104
34-DNT C	NA		#VALUE!	

**MS/MSD RECOVERY and RELATIVE PERCENT DIFFERENCE**

Non-Client

**LABORATORY CONTROL SAMPLES**

SAMPLE	COMPOUND	LCS %R	LCSD %R	RPD	ACTION
LFW791	12 of 17 HI		NA	NA	none, ND

Were laboratory acceptance limits used as the basis for validation? Yes  
 Did the laboratory provide CLP Form III or equivalent? Yes  
 Were chromatograms and quan reports present for all LCS/LCDs? Yes

LFW791

RDX

12 of 17 HI	AMOUNT FOUND	AMOUNT SPIKED	% R	FORM 3
LCS	1.27	1	127.0	127

%R = (Amount Found/Amount Spiked)\* 100

**BLANKS**

BLANK	COMPOUND	RESULT	5X OR 10X	ACTION LEVEL	ACTION
LFW791	135-TNBENZ	0.038	5	0.19	U at RL

ALL

NOTE Equipment/Field blanks are not qualified on the basis of laboratory method blank contamination or contamination in other field quality control blanks.

**CALIBRATION**

**Initial**

Are chromatograms and quan reports present for all ICAL standards? Yes  
 Are CLP For VIs or equivalent present and complete? Yes  
 Any transcription or calculation errors? No  
 What linearity criteria was used? 20% RSD or 0.99  
 Were RT window documented? Yes

**Continuing**

Are chromatograms and quan reports present for all CCV standards? Yes  
 Are CLP Form VIIs or equivalent present and complete? Yes  
 Any transcription or calculation errors? No  
 What %D criteria was used? 25% D  
 Have all samples and standards been listed on an Analytical Sequence? Yes  
 Was a proper analytical sequence followed? Yes

**INITIAL CALIBRATION - A**

Date	6/10/2009	Rptd Avg CF	77.89163	Rptd %RSD	5.156
Instrument	A	Calcd Avg CF	77.892	Calcd %RSD	5.156
Compound	RDX				
Reported CF	80.13	CF1	81.80	STD DEV	4
Calculated CF	80.42	CF2	75.90		
		CF3	77.15		
Response	16084	CF4	80.72		
Conc	200	CF5	78.45		
		CF6	80.13		
	CF = (H/C)	CF7	69.168	%RSD = (Std Dev/Avg CF)*100	
		CF8	79.815		

**CONTINUING CALIBRATION A**

Date	7/2/2009		
Time	1709		
Instrument	A		
Compound	HMX		
Reported Conc	104.5	Rptd %D	5
Calculated Conc	210.182	Calcd %D	-5.09
Response	25432	ICAL CF	200
CF	121	CCV CF	210.2
	CF = (H/C)	%Difference = ((ICAL - CCV)/ICAL)*100	

**FIELD DUPLICATES**

COMPOUND	SAMPLE	QUALIFIER	DUPLICATE	QUALIFIER	RPD
PCA					#DIV/0!

NOTES Samples are not qualified on this basis.

**COMPOUND IDENTIFICATION AND QUANTITATION**

Has a F10 been completed for every sample containing positive results? Yes  
 Was RT data presented on the form? NA  
 Are RTs within the established windows? Yes  
 Any transcription or calculation errors? No  
 Any false positives, negative peaks, shouldering, etc.? No  
 Was GCMS confirmation needed for results > 10 µg/ml? NA  
 Were percent differences or relative percent differences calculated? NA  
 Are percent differences/RPDs greater than 25%? NA  
 Are there any transcription errors? No  
 Are Form Is present for all field and quality control samples? No  
 Are chromatograms and quan reports present for all samples? Yes  
 Are RLs adjusted to reflect sample dilutions, percent solids, etc.? Yes  
 For soils, any percent solids <50%? No  
 For soils, any percent solids <10%? No

LCS RDX 1.27 µg/L

	A
response	4962
cf	77.89163
final vol ml	20
initial L	1
df	1
calculated	1.2741

23LC MW040W

TestAmerica West Sacramento

A-000006.D

Chromatography Summary

Injection Date: 7/2/2009 19:40 Operator: fhk  
 DataFile: LC10.I07022009.BVA-000006.D Vial Num: 13  
 Instrument ID: LC10

Method 8330 Target Analyte Results

Sample: **LFQAK1AC 9182192 G9F260326-1 1X** Method File: LC10.I07022009.B18330AB.M  
 Start Cal Date: 6/10/2009 15:33 End Cal Date: 6/11/2009 13:47

Matrix: WATER SubList: WATER sub SpikeList:  
 Samp. Info: LFQAK1AC 9182192 G9F260326-1 1X,0,  
 Misc. Info: ;,997 78.,20.1,WATER sub,;0;1,LFQAK1AC

Dilution Factor	Extract Volume	Sample Volume	Sample Weight
1X	20 mL	997.78 mL	0 g

Signal 1 UV 250-265						Signal 2 UV 358-205							
Compound Name	RT	Diff	Response	Conc (ug/L)	Flag	RT	Diff	Response	Conc (ug/L)	Flag	MDL	RL	Flag
3,4-Dinitrotoluene	18.24	0.001	6135	2.6070<		18.24	-0.006	11892	2.7570		0.0000	0.00	
HMX											0.0271	0.10	
RDX											0.0651	0.10	
Picric ACID											0.1253	1.00	
1,3,5-Trinitrobenzene	10.14	-0.086	264	0.0359<							0.0311	0.10	45
1,3-Dinitrobenzene											0.0501	0.10	
TETRYL											0.0501	0.10	
Nitrobenzene											0.0501	0.10	
2,4,6-Trinitrotoluene											0.0241	0.10	
4-AM-2,6-DNF											0.0220	0.10	
2-AM-4,6-DNF											0.1002	0.20	
2,6-Dinitrotoluene											0.0501	0.10	
2,4-Dinitrotoluene											0.0501	0.10	
2-Nitrotoluene											0.0722	0.50	
4-Nitrotoluene											0.0722	0.50	
3-Nitrotoluene											0.0621	0.50	
Nitroglycerin											0.3307	0.65	
PETN											0.3007	0.65	
3,5-Dinitroaniline											0.0251	1.00	

Surrogates:	Spiked	Recovered	% Rec	Spiked	Recovered	% Rec	Limits
3,4-Dinitrotoluene	2.5056	2.6070	104	2.5056	2.7570	110	(79-111)

Notes: M = Manually Integrated Signals Differ by More Than 40%  
 D = Operator Disabled Result Signals Differ by More Than 50%  
 O = Over Calibration Range

MATRIX SPIKE SAMPLE EVALUATION REPORT

HPLC

*Non-constant*

Client Lot #...: G9F260326      Work Order #...: LFV9K1AF-MS      Matrix.....: WATER  
 MS Lot-Sample #: G9F300242-004      LFV9K1AG-MSD  
 Date Sampled...: 06/26/09      Date Received...: 06/30/09  
 Prep Date.....: 07/01/09      Analysis Date...: 07/03/09  
 Prep Batch #...: 9182192  
 Dilution Factor: 0.98

PARAMETER	PERCENT RECOVERY	RECOVERY LIMITS	RPD	RPD LIMITS	METHOD
Nitroglycerin	106	(85 - 115)			SW846 8330
	108	(85 - 115)	2.0	(0-15)	SW846 8330
PETN	110	(84 - 117)			SW846 8330
	111	(84 - 117)	0.90	(0-15)	SW846 8330
Picric Acid	70	(21 - 118)			SW846 8330
	99 p	(21 - 118)	35	(0-20)	SW846 8330
2-Amino-4,6-dinitrotoluene	125 a	(77 - 123)			SW846 8330
	129 a	(77 - 123)	3.4	(0-27)	SW846 8330
4-Amino-2,6-dinitrotoluene	116 a	(68 - 113)			SW846 8330
	120 a	(68 - 113)	3.0	(0-30)	SW846 8330
1,3-Dinitrobenzene	125 a	(72 - 123)			SW846 8330
	130 a	(72 - 123)	4.1	(0-29)	SW846 8330
2,4-Dinitrotoluene	120 a	(70 - 119)			SW846 8330
	124 a	(70 - 119)	2.9	(0-30)	SW846 8330
2,6-Dinitrotoluene	122 a	(71 - 119)			SW846 8330
	125 a	(71 - 119)	2.9	(0-29)	SW846 8330
HMX	118 a	(67 - 115)			SW846 8330
	120 a	(67 - 115)	1.4	(0-32)	SW846 8330
Nitrobenzene	121 a	(69 - 119)			SW846 8330
	124 a	(69 - 119)	2.5	(0-31)	SW846 8330
2-Nitrotoluene	117	(64 - 120)			SW846 8330
	113	(64 - 120)	2.7	(0-36)	SW846 8330
3-Nitrotoluene	113	(67 - 114)			SW846 8330
	115 a	(67 - 114)	2.0	(0-31)	SW846 8330
4-Nitrotoluene	114	(67 - 115)			SW846 8330
	116 a	(67 - 115)	2.3	(0-32)	SW846 8330
RDX	121	(68 - 122)			SW846 8330
	129 a	(68 - 122)	6.3	(0-32)	SW846 8330
Tetryl	102	(66 - 105)			SW846 8330
	106 a	(66 - 105)	4.2	(0-26)	SW846 8330
1,3,5-Trinitrobenzene	119	(74 - 120)			SW846 8330
	124 a	(74 - 120)	4.3	(0-29)	SW846 8330
2,4,6-Trinitrotoluene	113 a	(69 - 111)			SW846 8330
	117 a	(69 - 111)	3.4	(0-28)	SW846 8330

(Continued on next page)

MATRIX SPIKE SAMPLE EVALUATION REPORT

HPLC

Client Lot #...: G9F260326      Work Order #...: LFV9K1AF-MS      Matrix.....: WATER  
MS Lot-Sample #: G9F300242-004      LFV9K1AG-MSD

<u>SURROGATE</u>	<u>PERCENT</u>	<u>RECOVERY</u>
	<u>RECOVERY</u>	<u>LIMITS</u>
3,4-Dinitrotoluene	104	(79 - 111)
	107	(79 - 111)

NOTE(S) :

Calculations are performed before rounding to avoid round-off errors in calculated results.

*Bold print denotes control parameters*

**p** Relative percent difference (RPD) is outside stated control limits

**a** Spiked analyte recovery is outside stated control limits



MATRIX SPIKE SAMPLE DATA REPORT

HPLC

Client Lot #...: G9F260326      Work Order #...: LFV9K1AF-MS      Matrix.....: WATER  
 MS Lot-Sample #: G9F300242-004      LFV9K1AG-MSD  
 Date Sampled...: 06/26/09      Date Received...: 06/30/09  
 Prep Date.....: 07/01/09      Analysis Date...: 07/03/09  
 Prep Batch #...: 9182192  
 Dilution Factor: 0.98

PARAMETER	SAMPLE	SPIKE	MEASRD	UNITS	PERCNT		METHOD
	AMOUNT	AMT	AMOUNT		RECVRY	RPD	
Nitroglycerin	ND	4.91	5.19	ug/L	106		SW846 8330
	ND	4.92	5.29	ug/L	108	2.0	SW846 8330
PETN	ND	4.91	5.41	ug/L	110		SW846 8330
	ND	4.92	5.46	ug/L	111	0.90	SW846 8330
Picric Acid	ND	4.91	3.41	ug/L	70		SW846 8330
	ND	4.92	4.85	ug/L	99 p	35	SW846 8330
2-Amino-4,6-dinitrotoluene	ND	0.982	1.22	ug/L	125 a		SW846 8330
	ND	0.983	1.26	ug/L	129 a	3.4	SW846 8330
4-Amino-2,6-dinitrotoluene	ND	0.982	1.14	ug/L	116 a		SW846 8330
	ND	0.983	1.18	ug/L	120 a	3.0	SW846 8330
1,3-Dinitrobenzene	ND	0.982	1.23	ug/L	125 a		SW846 8330
	ND	0.983	1.28	ug/L	130 a	4.1	SW846 8330
2,4-Dinitrotoluene	ND	0.982	1.18	ug/L	120 a		SW846 8330
	ND	0.983	1.22	ug/L	124 a	2.9	SW846 8330
2,6-Dinitrotoluene	ND	0.982	1.19	ug/L	122 a		SW846 8330
	ND	0.983	1.23	ug/L	125 a	2.9	SW846 8330
HMX	ND	0.982	1.16	ug/L	118 a		SW846 8330
	ND	0.983	1.18	ug/L	120 a	1.4	SW846 8330
Nitrobenzene	ND	0.982	1.19	ug/L	121 a		SW846 8330
	ND	0.983	1.22	ug/L	124 a	2.5	SW846 8330
2-Nitrotoluene	ND	0.982	1.14	ug/L	117		SW846 8330
	ND	0.983	1.11	ug/L	113	2.7	SW846 8330
3-Nitrotoluene	ND	0.982	1.11	ug/L	113		SW846 8330
	ND	0.983	1.14	ug/L	115 a	2.0	SW846 8330
4-Nitrotoluene	ND	0.982	1.12	ug/L	114		SW846 8330
	ND	0.983	1.14	ug/L	116 a	2.3	SW846 8330
RDX	ND	0.982	1.19	ug/L	121		SW846 8330
	ND	0.983	1.27	ug/L	129 a	6.3	SW846 8330
Tetryl	ND	0.982	0.998	ug/L	102		SW846 8330
	ND	0.983	1.04	ug/L	106 a	4.2	SW846 8330
1,3,5-Trinitrobenzene	0.035	0.982	1.20	ug/L	119		SW846 8330
	0.035	0.983	1.26	ug/L	124 a	4.3	SW846 8330
2,4,6-Trinitrotoluene	ND	0.982	1.11	ug/L	113 a		SW846 8330
	ND	0.983	1.15	ug/L	117 a	3.4	SW846 8330

(Continued on next page)

MATRIX SPIKE SAMPLE DATA REPORT

HPLC

Client Lot #...: G9F260326      Work Order #...: LFV9K1AF-MS      Matrix.....: WATER  
MS Lot-Sample #: G9F300242-004      LFV9K1AG-MSD

<u>SURROGATE</u>	<u>PERCENT RECOVERY</u>	<u>RECOVERY LIMITS</u>
3,4-Dinitrotoluene	104	(79 - 111)
	107	(79 - 111)

**NOTE(S) :**

Calculations are performed before rounding to avoid round-off errors in calculated results.

Bold print denotes control parameters

p Relative percent difference (RPD) is outside stated control limits.

a Spiked analyte recovery is outside stated control limits.

LABORATORY CONTROL SAMPLE EVALUATION REPORT

HPLC

Client Lot #...: G9F260326      Work Order #...: LFW791AC      Matrix.....: WATER  
 LCS Lot-Sample#: G9G010000-192  
 Prep Date.....: 07/01/09      Analysis Date...: 07/02/09  
 Prep Batch #...: 9182192  
 Dilution Factor: 1

<u>PARAMETER</u>	<u>PERCENT RECOVERY</u>	<u>RECOVERY LIMITS</u>	<u>METHOD</u>
Nitroglycerin	107	(85 - 115)	SW846 8330
PNTN	110	(84 - 117)	SW846 8330
Picric Acid	88	(21 - 118)	SW846 8330
2-Amino-4,6-dinitrotoluene	127 a	(77 - 123)	SW846 8330
4-Amino-2,6-dinitrotoluene	119 a	(68 - 113)	SW846 8330
1,3-Dinitrobenzene	128 a	(72 - 123)	SW846 8330
2,4-Dinitrotoluene	123 a	(70 - 119)	SW846 8330
2,6-Dinitrotoluene	125 a	(71 - 119)	SW846 8330
HMX	120 a	(67 - 115)	SW846 8330
Nitrobenzene	124 a	(69 - 119)	SW846 8330
2-Nitrotoluene	115	(64 - 120)	SW846 8330
3-Nitrotoluene	118 a	(67 - 114)	SW846 8330
4-Nitrotoluene	117 a	(67 - 115)	SW846 8330
RDX	127 a	(68 - 122)	SW846 8330
Tetryl	105	(66 - 105)	SW846 8330
1,3,5-Trinitrobenzene	127 a	(74 - 120)	SW846 8330
2,4,6-Trinitrotoluene	116 a	(69 - 111)	SW846 8330

<u>SURROGATE</u>	<u>PERCENT RECOVERY</u>	<u>RECOVERY LIMITS</u>
3,4-Dinitrotoluene	106	(79 - 111)

**NOTE(S) :**

Calculations are performed before rounding to avoid round-off errors in calculated results.

Bold print denotes control parameters

a Spiked analyte recovery is outside stated control limits.

LABORATORY CONTROL SAMPLE DATA REPORT

HPLC

Client Lot #...: G9F260326      Work Order #...: LFW791AC      Matrix.....: WATER  
 LCS Lot-Sample#: G9G010000-192  
 Prep Date.....: 07/01/09      Analysis Date...: 07/02/09  
 Prep Batch #...: 9182192  
 Dilution Factor: 1

<u>PARAMETER</u>	<u>SPIKE</u> <u>AMOUNT</u>	<u>MEASURED</u> <u>AMOUNT</u>	<u>UNITS</u>	<u>PERCENT</u> <u>RECOVERY</u>	<u>METHOD</u>
Nitroglycerin	5.00	5.35	ug/L	107	SW846 8330
PETN	5.00	5.52	ug/L	110	SW846 8330
Picric Acid	5.00	4.42	ug/L	88	SW846 8330
2-Amino-4,6-dinitrotoluene	1.00	1.27 a	ug/L	127	SW846 8330
4-Amino-2,6-dinitrotoluene	1.00	1.19 a	ug/L	119	SW846 8330
1,3-Dinitrobenzene	1.00	1.28 a	ug/L	128	SW846 8330
2,4-Dinitrotoluene	1.00	1.23 a	ug/L	123	SW846 8330
2,6-Dinitrotoluene	1.00	1.25 a	ug/L	125	SW846 8330
HMX	1.00	1.20 a	ug/L	120	SW846 8330
Nitrobenzene	1.00	1.24 a	ug/L	124	SW846 8330
2-Nitrotoluene	1.00	1.15	ug/L	115	SW846 8330
3-Nitrotoluene	1.00	1.18 a	ug/L	118	SW846 8330
4-Nitrotoluene	1.00	1.17 a	ug/L	117	SW846 8330
ROX	1.00	1.27 a	ug/L	127	SW846 8330
Tetryl	1.00	1.05	ug/L	105	SW846 8330
1,3,5-Trinitrobenzene	1.00	1.27 a	ug/L	127	SW846 8330
2,4,6-Trinitrotoluene	1.00	1.16 a	ug/L	116	SW846 8330

<u>SURROGATE</u>	<u>PERCENT</u> <u>RECOVERY</u>	<u>RECOVERY</u> <u>LIMITS</u>
3,4-Dinitrotoluene	106	(79 - 111)

**NOTE(S) :**

Calculations are performed before rounding to avoid round-off errors in calculated results.

Bold print denotes control parameters

a Spiked analyte recovery is outside stated control limits.

Chromatography Summary

Injection Date: 7/2/2009 18:50 Operator: fhk  
 DataFile: LC10 I\07022009 BVA-000005.D Vial Num: 12  
 Instrument ID: LC10

Method 8330 Target Analyte Results

Sample : LFW791AC 9182192 G9G010000-192  
 C 1X

Method File: LC10 I\07022009 B\8330AB.M  
 Start Cal Date: 6/10/2009 15:33 End Cal Date: 6/11/2009 13:47

Matrix: WATER SubList: WATER sub SpikeList: WATER spk  
 Samp. Info: LFW791AC 9182192 G9G010000-192 C 1X:3;  
 Misc. Info: LCS.,1000.,20;1-WATER sub;WATER spk,1,1,LFW791AC

Dilution Factor	Extract Volume	Sample Volume	Sample Weight
1X	20 mL	1000 mL	0 g

Signal 1 UV 250-265								Signal 2 UV 358-205								
Compound Name	RT	Response	Conc (ug/L)	Spike Level	%R	Result	Flag	RT	Response	Conc (ug/L)	Spike Level	%R	Result	Flag	Limits(%)	Flag
3,4-Dinitrotoluene	18.23	6246	2.6480<	2.5	106%	Acceptable		18.23	12082	2.7950	2.5	112%	Fails		(79-111)	
HMX	5.30	7265	1.2020<	1	120%	Fails					1	0%	Fails		(65-115)	45
RDX	7.88	4962	1.2740<	1	127%	Fails					1	0%	Fails		(68-122)	45
Picric ACID	9.06	18341	4.4320	5	89%	Acceptable		9.06	26909	4.4230<	5	88%	Acceptable		(21-118)	
1,3,5-Trinitrobenzene	10.23	9382	1.2710<	1	127%	Fails					1	0%	Fails		(74-120)	45
1,3-Dinitrobenzene	13.30	8977	1.2850<	1	129%	Fails					1	0%	Fails		(72-123)	45
TETRYL	14.35	4859	1.0470<	1	105%	Acceptable					1	0%	Fails		(66-105)	45
Nitrobenzene	15.23	3778	1.2450<	1	125%	Fails					1	0%	Fails		(69-119)	45
2,4,6-Trinitrotoluene	16.58	5089	1.1600<	1	116%	Fails					1	0%	Fails		(69-111)	45
4-AM-2,6-DNT	17.60	3716	1.1930<	1	119%	Fails					1	0%	Fails		(68-113)	45
2-AM-4,6-DNT	18.66	4662	1.2680<	1	127%	Fails					1	0%	Fails		(77-123)	45
2,6-Dinitrotoluene	20.39	3326	1.2490<	1	125%	Fails					1	0%	Fails		(71-119)	45
2,4-Dinitrotoluene	21.10	5280	1.2330<	1	123%	Fails					1	0%	Fails		(70-119)	45
2-Nitrotoluene	24.82	2061	1.1470<	1	115%	Acceptable					1	0%	Fails		(64-120)	45
4-Nitrotoluene	26.68	2553	1.1660<	1	117%	Fails					1	0%	Fails		(67-115)	45
3-Nitrotoluene	28.69	2534	1.1750<	1	118%	Fails					1	0%	Fails		(67-114)	45
Nitroglycerin				5	0%	Fails		15.84	16262	5.3460<	5	107%	Acceptable		(85-115)	45
PETN				5	0%	Fails		30.25	8838	5.5150<	5	110%	Acceptable		(84-117)	45
3,5-Dinitroaniline	14.04	5620	1.2490<	1	125%	Acceptable					1	0%	Fails		(40-140)	45

Surrogates	Spiked	Recovered	% Rec	Spiked	Recovered	% Rec	Limits
3,4-Dinitrotoluene	2.5000	2.6480	106	2.5000	2.7950	112	(79-111)

Notes: M = Manually Integrated Signals Differ by More Than 40%  
 D = Operator Disabled Result Signals Differ by More Than 50%  
 O = Over Calibration Range

LCS

Test America West Sacramento

Method 8330

Data file : \\Terastation\share\GCdata\LC10.I\07022009.B\A-000005.D  
 Lab Smp Id: LFW791AC 9182192 G9  
 Inj Date : 02-JUL-2009 18:50  
 Operator : fhk Inst ID: LC10.i  
 Smp Info : LFW791AC 9182192 G9G010000-192 C 1X;3;  
 Misc Info : LCS;;1000;;20;1;WATER.sub;WATER.spk;1;1;LFW791AC  
 Comment : SOP SAC-LC-0009  
 Method : \\Terastation\share\GCdata\LC10.I\07022009.B\8330AB.m  
 Meth Date : 02-Jul-2009 19:27 kenneyf Quant Type: AREA%  
 Cal Date : 11-JUN-2009 13:47 Cal File: A-000015.d  
 Als bottle: 12 QC Sample: LCS  
 Dil Factor: 1.00000  
 Integrator: Falcon+ Compound Sublist: WATER.sub  
 Target Version: 4.14  
 Processing Host: SACP307HPLC

Concentration Formula: Amt \* DF \* Uf\*Vt / Vo \* CpndVariable

Name	Value	Description
DF	1.000	Dilution Factor
Uf	1.000	ng unit correction factor
Vt	20.000	Volume of final extract (mL)
Vo	1000.000	Volume of sample extracted (mL)
Cpnd Variable		Local Compound Variable

RT	AREA	HEIGHT	HT/AREA	% HEIGHT	COMPOUNDS
5.298	44338	7265	0.164	7.31	2 HMX
7.878	52369	4962	0.095	4.99	3 RDX
9.061	240427	18341	0.076	18.58	5 Picric ACID
10.228	121534	9382	0.077	9.44	6 1,3,5-Trinitrobenz
11.635	533	49	0.092	0.04	
12.105	7464	514	0.069	0.51	
12.778	132	46	0.347	0.04	
13.301	143585	8977	0.063	9.04	7 1,3-Dinitrobenzene
14.045	89298	5620	0.063	5.66	8 3,5-Dinitroaniline
14.351	82373	4859	0.059	4.89	9 TETRYL
15.235	67423	3778	0.056	3.80	10 Nitrobenzene
16.585	95229	5089	0.053	5.12	12 2,4,6-Trinitrotolu
17.598	75159	3716	0.049	3.74	13 4-AM-2,6-DNT
18.235	122827	6246	0.051	6.29	1 3,4-Dinitrotoluene
18.658	104810	4662	0.044	4.69	14 2-AM-4,6-DNT
20.395	74253	3326	0.045	3.34	15 2,6-Dinitrotoluene
21.105	128415	5280	0.041	5.31	16 2,4-Dinitrotoluene
23.125	284	28	0.099	0.02	
24.825	56586	2061	0.036	2.07	17 2-Nitrotoluene
26.678	75508	2553	0.034	2.57	18 4-Nitrotoluene
28.688	80096	2534	0.032	2.55	19 3-Nitrotoluene
	1662642	99288		100.000	

Total unknown % height = 0.6100



METHOD BLANK REPORT

HPLC

Client Lot #...: G9F260326      Work Order #...: LFW791AA      Matrix.....: WATER  
 MB Lot-Sample #: G9G010000-192  
 Analysis Date...: 07/02/09      Prep Date.....: 07/01/09  
 Dilution Factor: 1      Prep Batch #...: 9182192

PARAMETER	RESULT	REPORTING LIMIT	UNITS	METHOD
Nitroglycerin	ND	0.65	ug/L	SW846 8330
PETN	ND	0.65	ug/L	SW846 8330
Picric Acid	ND	1.0	ug/L	SW846 8330
2-Amino-4,6-dinitrotoluene	ND	0.20	ug/L	SW846 8330
4-Amino-2,6-dinitrotoluene	ND	0.10	ug/L	SW846 8330
1,3-Dinitrobenzene	ND	0.10	ug/L	SW846 8330
2,4-Dinitrotoluene	ND	0.10	ug/L	SW846 8330
2,6-Dinitrotoluene	ND	0.10	ug/L	SW846 8330
HMX	ND	0.10	ug/L	SW846 8330
Nitrobenzene	ND	0.10	ug/L	SW846 8330
2-Nitrotoluene	ND	0.40	ug/L	SW846 8330
3-Nitrotoluene	ND	0.40	ug/L	SW846 8330
4-Nitrotoluene	ND	0.50	ug/L	SW846 8330
RDX	ND	0.10	ug/L	SW846 8330
Tetryl	ND	0.10	ug/L	SW846 8330
1,3,5-Trinitrobenzene	0.039 J	0.10	ug/L	SW846 8330
2,4,6-Trinitrotoluene	ND	0.10	ug/L	SW846 8330

SURROGATE	PERCENT RECOVERY	RECOVERY LIMITS
3,4-Dinitrotoluene	106	(79 - 111)

NOTE(S):

Calculations are performed before rounding to avoid round-off errors in calculated results.

J Estimated result. Result is less than RL

Data File: \\Terastation\share\GCdata\LC10.I\07022009.B\A-000004.D

Date : 02-JUL-2009 18:00

Instrument: LC10.i

Client ID:

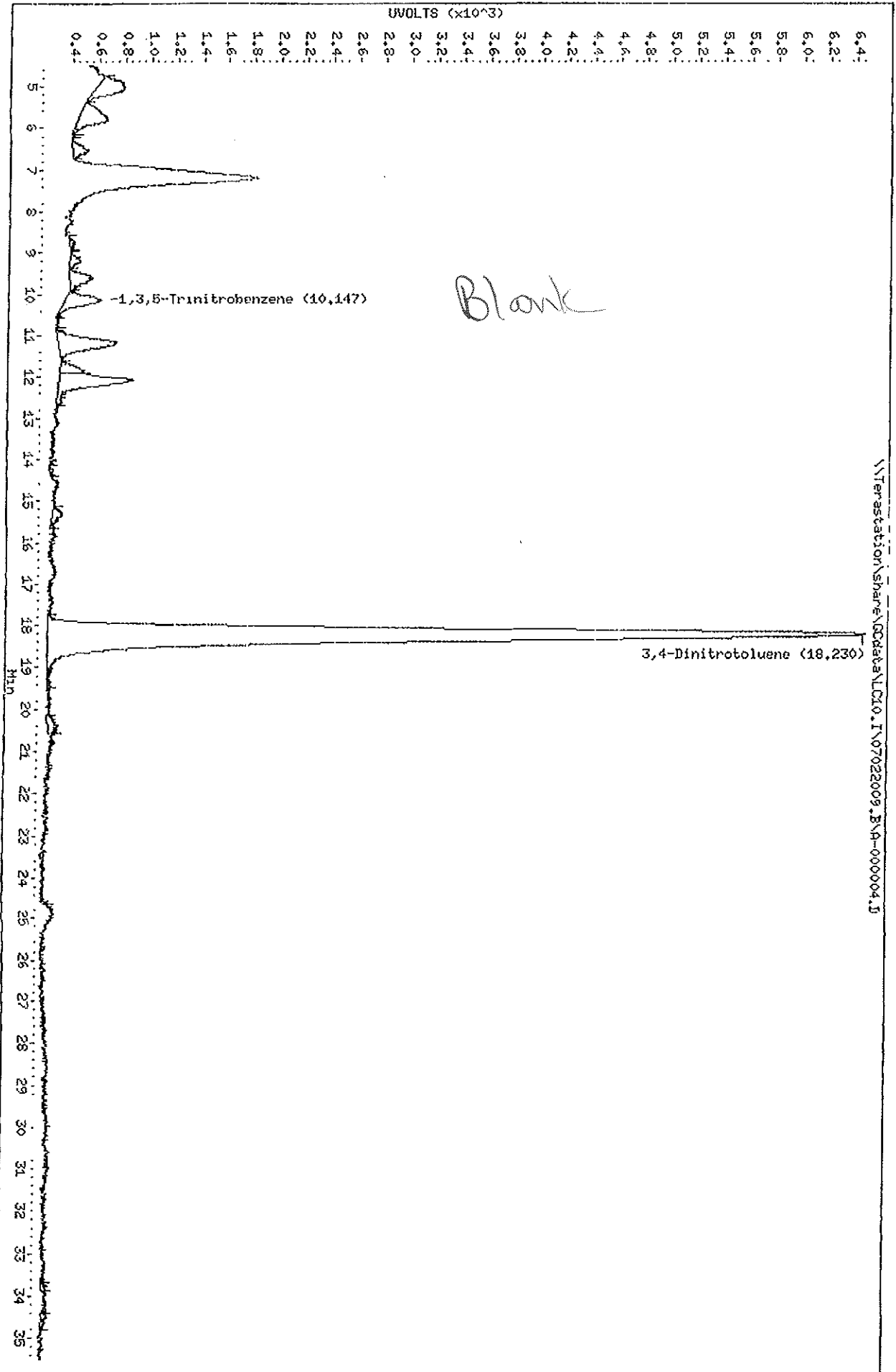
Sample Info: LFM794AA 9182192 GPC010000-192 B 1X:09

Volume Injected (uL): 500.0

Column phase: SYNERGI HYDRORP C18

Operator: PHK  
Column diameter: 4.60

\\Terastation\share\GCdata\LC10.I\07022009.B\A-000004.D



# QC DATA ASSOCIATION SUMMARY

G9F260326

Sample Preparation and Analysis Control Numbers

<u>SAMPLE#</u>	<u>MATRIX</u>	<u>ANALYTICAL METHOD</u>	<u>LEACH BATCH #</u>	<u>PREP BATCH #</u>	<u>MS RUN#</u>
001	WATER	SW846 8330		9182192	9182194
002	WATER	SW846 8330		9182192	9182194
003	WATER	SW846 8330		9182192	9182194



RQC058

TestAmerica Laboratories, Inc.  
EXTRACTION BENCH WORKSHEET

Run Date: 7/02/09  
Time: 11:49:39

LEV 1	LEV 1	Blank	Weights/Volumes
Y 2	Y 2	Check	Spike & Surrogate Worksheet
Y 1	Y 1	MS/MSD	Vial contains correct volume
			Labels, greenbars, worksheets
			computer batch: correct & all match
			Anomalies to Extraction Method

Expanded Deliverable  
COC Completed  
Y Bench Sheet Copied  
Package Submitted to Analytical Group  
Bench Sheet Copied per COC

Extractionist: 002448 Tuan Q. Phan

PREP DATE: 7/01/09 9:00  
COMP DATE: 7/02/09 10:40

\*\*\*\*\*  
\* QC BATCH: 9182192 \*  
\* \*\*\*\*\* \*

Reviewer/Date: PHANT / 7/02/09

Nitroaromatics & Nitramines: Explosives (8330)  
SOLID PHASE EXTRACTION (NOMINAL)

EXTR EXPR	ANL DUE	LOT# MSRUN# / WORK ORDER	TEST FLGS	EXT MTH	MATRIX	INIT/VOL	PH#S ADJ1	INIT ADJ2	EXTRACTION VOL	SOLVENTS VOL EXCHANGE	VOL	SPIKE STANDARD / SURROGATE ID	
7/01/09	7/10/09	G9F260326-001 LFOAK-1-AC		B7	A0 WATER	997.78mL 20.00mL	NA	NA	HOAC/ACN	4.5	HOAC/ACN	5.0	50UL-09GCSV0172
COMMENTS:													
7/01/09	7/10/09	G9F260326-002 LFOAL-1-AC		B7	A0 WATER	991.12mL 20.00mL	NA	NA	HOAC/ACN	4.5	HOAC/ACN	5.0	50UL-09GCSV0172
COMMENTS:													
7/01/09	7/10/09	G9F260326-003 LFOAM-1-AC		B7	A0 WATER	1009.82mL 20.00mL	NA	NA	HOAC/ACN	4.5	HOAC/ACN	5.0	50UL-09GCSV0172
COMMENTS:													
7/01/09	7/13/09	G9F270189-001 LFRD6-1-AC		B7	A0 WATER	995.56mL 20.00mL	NA	NA	HOAC/ACN	4.5	HOAC/ACN	5.0	50UL-09GCSV0172
COMMENTS:													
7/02/09	7/13/09	G9F270189-002 LFRD7-1-AC		B7	A0 WATER	999.12mL 20.00mL	NA	NA	HOAC/ACN	4.5	HOAC/ACN	5.0	50UL-09GCSV0172
COMMENTS:													
7/02/09	7/13/09	G9F270189-003 LFRD8-1-AC		B7	A0 WATER	991.73mL 20.00mL	NA	NA	HOAC/ACN	4.5	HOAC/ACN	5.0	50UL-09GCSV0172
COMMENTS:													
7/02/09	7/13/09	G9F270189-004 LFRD9-1-AC		B7	A0 WATER	996.78mL 20.00mL	NA	NA	HOAC/ACN	4.5	HOAC/ACN	5.0	50UL-09GCSV0172
COMMENTS:													

RQC058

TestAmerica Laboratories, Inc.  
EXTRACTION BENCH WORKSHEET

Run Date: 7/02/09  
Time: 11:49:39

\*\*\*\*\*  
\* QC BATCH: 9182192 \*  
\* PREP DATE: 7/01/09 9:00  
\* COMP DATE: 7/02/09 10:40  
\*\*\*\*\*

EXTR EXPR	ANL DUE	LOT# WORK ORDER	MSRIN#/ ORDER	TEST FLGS	EXT MTH	MATRIX	INIT/FIN WT/VOL	PH'S ADJ1	ADJ2	EXTRACTION VOL	SOLVENTS VOL	EXCHANGE	VOL	SPIKE STANDARD/ SURROGATE ID
7/02/09	7/13/09	G9F270189-005	LFV9A-1-AC		B7	A0	985.33mL 20.00mL	NA	NA	HOAC/ACN	4.5	HOAC/ACN	5.0	50UL-09GCSV0172
COMMENTS:														
7/02/09	7/14/09	G9F300242-001	LFV9A-1-AC		B7	A0	1014.4mL 20.00mL	NA	NA	HOAC/ACN	4.5	HOAC/ACN	5.0	50UL-09GCSV0172
COMMENTS:														
7/02/09	7/14/09	G9F300242-002	LFV9G-1-AC		B7	A0	975.28mL 20.00mL	NA	NA	HOAC/ACN	4.5	HOAC/ACN	5.0	50UL-09GCSV0172
COMMENTS:														
7/02/09	7/14/09	G9F300242-003	LFV9J-1-AC		B7	A0	1018.21mL 20.00mL	NA	NA	HOAC/ACN	4.5	HOAC/ACN	5.0	50UL-09GCSV0172
COMMENTS:														
7/03/09	7/14/09	G9F300242-004	LFV9K-1-AC		B7	A0	1028.11mL 20.00mL	NA	NA	HOAC/ACN	4.5	HOAC/ACN	5.0	50UL-09GCSV0172
COMMENTS:														
7/03/09	7/14/09	G9F300242-004	LFV9K-1-AFS		B7	A0	1018.28mL 20.00mL	NA	NA	HOAC/ACN	4.5	HOAC/ACN	5.0	SEE BENCH SHEET 50UL-09GCSV0172
COMMENTS:														
7/03/09	7/14/09	G9F300242-004	LFV9K-1-AGD		B7	A0	1017.3mL 20.00mL	NA	NA	HOAC/ACN	4.5	HOAC/ACN	5.0	SEE BENCH SHEET 50UL-09GCSV0172
COMMENTS:														
7/03/09	7/14/09	G9F300242-005	LFV9M-1-AC		B7	A0	1016.72mL 20.00mL	NA	NA	HOAC/ACN	4.5	HOAC/ACN	5.0	50UL-09GCSV0172
COMMENTS:														
7/01/09	0/00/00	G9G010000-192	LFW79-1-AAB		B7	A0	1000mL 20.00mL	NA	NA	HOAC/ACN	4.5	HOAC/ACN	5.0	50UL-09GCSV0172
COMMENTS:														
7/01/09	0/00/00	G9G010000-192	LFW79-1-ACC		B7	A0	1000mL 20.00mL	NA	NA	HOAC/ACN	4.5	HOAC/ACN	5.0	SEE BENCH SHEET 50UL-09GCSV0172
COMMENTS:														

1% HOAC/ACN 2991-95F; 1%HOAC/H2O 2991-87B; FILTER MILLIPORE LOT R5FN29184  
 SPE COLUMN WATER LOT 00339084A; SODIUM CHLORIDE MALLINKRODT LOT H07582  
 LCS.MS/MSD SPIKE 20UL-09GCSV0037;100UL-09GCSV0149;100UL-09GCSV0039

TestAmerica West Sacramento  
GC/LC INSTRUMENT LOG

Inst ID: LC10                      Batch ID: 06102009  
Method : Method 8330              Test : SOP SAC-LC-0009  
ICAL Date: See Calibration Report

Date	Time	Operator	Sample ID	File ID	Vol or wt	Extract Vol	Diln	Comments
10-JUN-2009	14:42	fhk	Blank	A-000001.	0 g	0 mL	1	
10-JUN-2009	15:33	fhk	STD_1 09GCSV0048 5.0/0/0/0ng/ml	A-000002.	0 g	0 mL	1	
10-JUN-2009	16:23	fhk	STD_2 09GCSV0049 10/20/10/10ng	A-000003.	0 g	0 mL	1	
10-JUN-2009	17:14	fhk	STD_3 09GCSV0050 20/50/20/20ng	A-000004.	0 g	0 mL	1	Bad Std
10-JUN-2009	18:05	fhk	STD_4 09GCSV0051 50/100/50/50n	A-000005.	0 g	0 mL	1	
10-JUN-2009	18:56	fhk	STD_5 09GCSV0053 100/200/100/1	A-000006.	0 g	0 mL	1	
10-JUN-2009	19:46	fhk	STD_6 09GCSV0054 200/500/200/2	A-000007.	0 g	0 mL	1	
10-JUN-2009	20:37	fhk	STD_7 09GCSV0055 500/1000/500/	A-000008.	0 g	0 mL	1	
10-JUN-2009	21:28	fhk	STD_8 09GCSV0056 1000/2000/100	A-000009.	0 g	0 mL	1	
10-JUN-2009	22:18	fhk	Blank	A-000010.	0 g	0 mL	1	
10-JUN-2009	23:09	fhk	ICV_6 09GCSV0397 100/200/100/1	A-000011.	0 g	0 mL	1	
11-JUN-2009	00:00	fhk	STD_5 09GCSV0053 100/200/100/1	A-000012.	0 g	0 mL	1	
11-JUN-2009	00:50	fhk	Surrogate 100ng/mL	A-000013.	0 g	0 mL	1	
11-JUN-2009	12:56	fhk	Primer	A-000014.	0 g	0 mL	1	
11-JUN-2009	13:47	fhk	STD_3 09GCSV0050 20/50/20/20ng	A-000015.	0 g	0 mL	1	was injected Nov 5/09

*JM 6/11/09*



TestAmerica West Sacramento

INITIAL CALIBRATION DATA

Start Cal Date : 10-JUN-2009 15:33  
 End Cal Date : 11-JUN-2009 13:47  
 Quant Method : ESTD  
 Origin : Disabled  
 Target Version : 4.14  
 Integrator : Falcon  
 Method file : \\Terastation\share\GCdata\LC10.I\06102009.B\8330AB.M  
 Last Edit : 11-Jun-2009 15:06 kenneyf  
 Curve Type : Average

Calibration File Names:

Level 1: \\Terastation\share\GCdata\GCdata\LC10.I\06102009.B\A-000002.d  
 Level 2: \\Terastation\share\GCdata\GCdata\LC10.I\06102009.B\A-000003.d  
 Level 3: \\Terastation\share\GCdata\GCdata\LC10.I\06102009.B\A-000015.d  
 Level 4: \\Terastation\share\GCdata\GCdata\LC10.I\06102009.B\A-000005.d  
 Level 5: \\Terastation\share\GCdata\GCdata\LC10.I\06102009.B\A-000006.d  
 Level 6: \\Terastation\share\GCdata\GCdata\LC10.I\06102009.B\A-000007.d  
 Level 7: \\Terastation\share\GCdata\GCdata\LC10.I\06102009.B\A-000008.d  
 Level 8: \\Terastation\share\GCdata\GCdata\LC10.I\06102009.B\A-000009.d

Compound	5.000	10.000	20.000	50.000	100.000	200.000	RRF	% RSD
	Level 1	Level 2	Level 3	Level 4	Level 5	Level 6		
	500.000	1000.000						
	Level 7	Level 8						
2 HMX	119 ✓ 116	120 ✓ 123	115 ✓	125 ✓	123 ✓	124	121	3.066
3 RDX	81.80000 ✓ 69.16800	75.90000 ✓ 79.81500	77.15000 ✓	80.72000 ✓	78.45000 ✓	80.13000	77.89163	5.156
4 EGDN	++++ ++++	++++ ++++	++++	++++	++++	++++	++++	++++
5 Picric ACID	++++ 74.63400	++++ 74.87800	91.72000	85.66000	84.85500	84.88200	82.77483	8.120
6 1,3,5-Trinitrobenzene	161 136	146 147	145	150	147	148	148	4.610
7 1,3-Dinitrobenzene	145 128	142 141	138	142	139	142	140	3.715
8 3,5-Dinitroaniline	97.20000 82.20800	90.80000 89.25000	89.35000	91.80000	88.83000	90.35000	89.97350	4.581
9 TETRYL	97.60000 ✓ 92.43000	95.10000 ✓ 94.40900	77.40000 ✓	96.02000 ✓	94.31000 ✓	95.06000	92.79113	6.888

*AK*  
6/11/09

TestAmerica West Sacramento

INITIAL CALIBRATION DATA

Start Cal Date : 10-JUN-2009 15:33  
 End Cal Date : 11-JUN-2009 13:47  
 Quant Method : ESTD  
 Origin : Disabled  
 Target Version : 4.14  
 Integrator : Falcon  
 Method file : \\Terastation\share\GCdata\LC10.I\06102009.B\8330AB.M  
 Last Edit : 11-Jun-2009 15:06 kenneyf  
 Curve Type : Average

Compound	5.000	10.000	20.000	50.000	100.000	200.000	RRF	% RSD
	Level 1	Level 2	Level 3	Level 4	Level 5	Level 6		
	500.000	1000.000						
	Level 7	Level 8						
10 Nitrobenzene	63.80000 56.02600	62.30000 61.95700	56.55000	61.44000	61.06000	62.26000	60.67412	4.659
11 Nitroglycerin	+++++	+++++	+++++	+++++	+++++	+++++	+++++	+++++
12 2,4,6-Trinitrotoluene	94.80000 84.22200	90.20000 87.17800	91.90000	88.78000	87.04000	87.55500	87.70937	4.401
13 4-AM-2,6-DNT	67.00000 58.65400	63.30000 61.14600	60.85000	63.26000	61.61000	62.67000	62.31125	3.901
14 2-AM-4,6-DNT	74.80000 69.18400	75.80000 72.13500	73.10000	75.54000	73.53000	74.39000	73.55987	2.932
15 2,6-Dinitrotoluene	57.60000 50.43000	53.30000 52.74900	51.90000	53.68000	53.03000	53.21500	53.23800	3.835
16 2,4-Dinitrotoluene	93.00000 81.43200	86.50000 85.05700	82.60000	86.20000	84.72000	85.49000	85.62488	4.031
17 2-Nitrotoluene	40.80000 33.97400	37.40000 35.45000	33.20000	35.64000	35.21000	35.77500	35.93113	6.486
18 4-Nitrotoluene	48.00000 41.67600	46.40000 43.21300	40.75000	43.56000	43.22000	43.56500	43.79800	5.390
19 3-Nitrotoluene	47.00000 41.51600	44.60000 42.82400	39.85000	43.24000	42.93000	43.08000	43.13000	4.852
20 PETN	+++++	+++++	+++++	+++++	+++++	+++++	+++++	+++++

TestAmerica West Sacramento

INITIAL CALIBRATION DATA

Start Cal Date : 10-JUN-2009 15:33  
 End Cal Date : 11-JUN-2009 13:47  
 Quant Method : ESTD  
 Origin : Disabled  
 Target Version : 4.14  
 Integrator : Falcon  
 Method file : \\Terastation\share\GCdata\LC10.I\06102009.B\8330AB.M  
 Last Edit : 11-Jun-2009 15:06 kenneyf  
 Curve Type : Average

Compound	5.000 Level 1	10.000 Level 2	20.000 Level 3	50.000 Level 4	100.000 Level 5	200.000 Level 6	RRF	% RSD
	500.000 Level 7	1000.000 Level 8						
\$ 1 3,4-Dinitrotoluene	++++	50.10000	46.40000	46.48000	45.87000	46.85500	47.16729	3.251
	46.08000	48.38600						

TestAmerica West Sacramento

INITIAL CALIBRATION DATA

Start Cal Date : 10-JUN-2009 16:23  
 End Cal Date : 11-JUN-2009 13:47  
 Quant Method : ESTD  
 Origin : Disabled  
 Target Version : 4.14  
 Integrator : Falcon  
 Method file : \\Terastation\share\GCdata\LC10.I\06102009.B\8330AB.M\833  
 Last Edit : 11-Jun-2009 15:14 kenneyf  
 Curve Type : Average

Calibration File Names:

Level 2: \\Terastation\share\GCdata\GCdata\LC10.I\06102009.B\A-000003.d\A-C  
 Level 3: \\Terastation\share\GCdata\GCdata\LC10.I\06102009.B\A-000015.d\A-C  
 Level 4: \\Terastation\share\GCdata\GCdata\LC10.I\06102009.B\A-000005.d\A-C  
 Level 5: \\Terastation\share\GCdata\GCdata\LC10.I\06102009.B\A-000006.d\A-C  
 Level 6: \\Terastation\share\GCdata\GCdata\LC10.I\06102009.B\A-000007.d\A-C  
 Level 7: \\Terastation\share\GCdata\GCdata\LC10.I\06102009.B\A-000008.d\A-C  
 Level 8: \\Terastation\share\GCdata\GCdata\LC10.I\06102009.B\A-000009.d\A-C

Compound	10.000 Level 2	20.000 Level 3	50.000 Level 4	100.000 Level 5	200.000 Level 6	500.000 Level 7	Level 8	RRF	% RSD
2 HMX	+++++	+++++	+++++	+++++	+++++	+++++	+++++	+++++	+++++
3 RDX	+++++	+++++	+++++	+++++	+++++	+++++	+++++	+++++	+++++
4 EGDN	+++++	+++++	+++++	+++++	+++++	+++++	+++++	+++++	+++++
5 Picric ACID	+++++ 110	134	126	125	125	110		122	8.080
6 1,3,5-Trinitrobenzene	+++++	+++++	+++++	+++++	+++++	+++++	+++++	+++++	+++++
7 1,3-Dinitrobenzene	+++++	+++++	+++++	+++++	+++++	+++++	+++++	+++++	+++++
8 3,5-Dinitroaniline	+++++	+++++	+++++	+++++	+++++	+++++	+++++	+++++	+++++
9 TETRYL	+++++	+++++	+++++	+++++	+++++	+++++	+++++	+++++	+++++

TestAmerica West Sacramento

INITIAL CALIBRATION DATA

Start Cal Date : 10-JUN-2009 16:23  
 End Cal Date : 11-JUN-2009 13:47  
 Quant Method : ESTD  
 Origin : Disabled  
 Target Version : 4.14  
 Integrator : Falcon  
 Method file : \\Terastation\share\GCdata\LC10.I\06102009.B\8330AB.M\833  
 Last Edit : 11-Jun-2009 15:14 kenneyf  
 Curve Type : Average

Compound	10.000 Level 2	20.000 Level 3	50.000 Level 4	100.000 Level 5	200.000 Level 6	500.000 Level 7	RRF	% RSD
	1000.000 Level 8							
10 Nitrobenzene	++++	++++	++++	++++	++++	++++	++++	++++
11 Nitroglycerin	++++ 61.73900	61.25000	59.82000	60.58000	61.86500	59.75400	60.83467	1.527
12 2,4,6-Trinitrotoluene	++++	++++	++++	++++	++++	++++	++++	++++
13 4-AM-2,6-DNT	++++	++++	++++	++++	++++	++++	++++	++++
14 2-AM-4,6-DNT	++++	++++	++++	++++	++++	++++	++++	++++
15 2,6-Dinitrotoluene	++++	++++	++++	++++	++++	++++	++++	++++
16 2,4-Dinitrotoluene	++++	++++	++++	++++	++++	++++	++++	++++
17 2-Nitrotoluene	++++	++++	++++	++++	++++	++++	++++	++++
18 4-Nitrotoluene	++++	++++	++++	++++	++++	++++	++++	++++
19 3-Nitrotoluene	++++	++++	++++	++++	++++	++++	++++	++++
20 PETN	++++ 32.75200	32.65000	30.36000	31.91000	32.25000	32.38000	32.05033	2.748

TestAmerica West Sacramento

INITIAL CALIBRATION DATA

Start Cal Date : 10-JUN-2009 16:23  
 End Cal Date : 11-JUN-2009 13:47  
 Quant Method : ESTD  
 Origin : Disabled  
 Target Version : 4.14  
 Integrator : Falcon  
 Method file : \\Terastation\share\GCdata\LC10.I\06102009.B\8330AB.M\83:  
 Last Edit : 11-Jun-2009 15:14 kenneyf  
 Curve Type : Average

Compound	10.000 Level 2	20.000 Level 3	50.000 Level 4	100.000 Level 5	200.000 Level 6	500.000 Level 7	RRF	% RSD
	1000.000 Level 8							
\$ 1 3,4-Dinitrotoluene	* +++++ 93.14200	90.40000	88.76000	88.28000	90.51000	88.03667	89.85478	2.142

*\* Level 2 dropped due to poor integration  
 Jbx  
 6/12/09*

Chromatography Summary

Injection Date: 6/10/2009 23:09 Operator: fhk  
 Data File: LC10.N06102009.BVA-000011.D Vial Num: 20  
 Instrument ID: LC10

Method 8330 Target Analyte Results

Sample: **ICV\_6 08GCSV0397**  
**100/200/100/100ng/mL**

Method File: LC10.N06102009.BVA8330AB.M  
 Start Cal Date: 6/10/2009 15:33 End Cal Date: 6/11/2009 13:47

Matrix: NONE SubList: CAL.sub SpikeList:  
 Samp. Info: ICV\_6 08GCSV0397 100/200/100/100ng/mL.2  
 Misc. Info: ;0; ; ;3;CAL.sub: ;0;1

Dilution Factor	Extract Volume	Sample Volume	Sample Weight
1X	0 mL	0 mL	0 g

Compound Name	RT	Signal 1 UV 250-265					Flag	Signal 2 UV 358-205					Limits(%)	Flag
		Response	PPB	Spike Level	%D	Result		RT	Response	PPB	Spike Level	%D		
<b>3,4-Dinitrotoluene</b>				200	-100%	Fails				200	-100%	Fails	(±15)	
HMX	5.29	26153	216.4000<	200	3%	Acceptable				200	-100%	Fails	(±15)	45
RDX	7.84	16084	206.5000<	200	3%	Acceptable				200	-100%	Fails	(±15)	45
Picric ACID	9.00	41754	504.4000 ✓	500	1%	Acceptable	9.00	61509	505.5000<	500	1%	Acceptable	(±15)	
1,3,5-Trinitrobenzene	10.20	29097	197.1000<	200	-1%	Acceptable				200	-100%	Fails	(±15)	45
1,3-Dinitrobenzene	13.25	27834	199.2000<	200	0%	Acceptable				200	-100%	Fails	(±15)	45
TBTRYL	14.30	17176	191.6000<	200	-4%	Acceptable				200	-100%	Fails	(±15)	45
Nitrobenzene	15.16	12143	200.1000<	200	0%	Acceptable				200	-100%	Fails	(±15)	45
2,4,6-Trinitrotoluene	16.53	17352	197.8000<	200	-1%	Acceptable				200	-100%	Fails	(±15)	45
4-AM-2,6-DNT	17.51	11969	192.1000< ✓	200	-4%	Acceptable				200	-100%	Fails	(±15)	45
2-AM-4,6-DNT	18.57	13851	188.3000<	200	-6%	Acceptable				200	-100%	Fails	(±15)	45
2,6-Dinitrotoluene	20.28	10445	196.2000<	200	-2%	Acceptable				200	-100%	Fails	(±15)	45
2,4-Dinitrotoluene	21.00	16653	194.5000<	200	-3%	Acceptable				200	-100%	Fails	(±15)	45
2-Nitrotoluene	24.68	7069	196.7000<	200	-2%	Acceptable				200	-100%	Fails	(±15)	45
4-Nitrotoluene	26.50	8526	194.7000<	200	-3%	Acceptable				200	-100%	Fails	(±15)	45
3-Nitrotoluene	28.49	8451	195.9000< ✓	200	-2%	Acceptable				200	-100%	Fails	(±15)	45
<b>Nitroglycerin</b>				200	-100%	Fails	15.78	12548	206.3000<	200	3%	Acceptable	(±15)	45
PETN				200	-100%	Fails	30.10	5830	181.9000< ✓	200	-9%	Acceptable	(±15)	45
3,5-Dinitroaniline	13.99	18460	205.2000<	200	3%	Acceptable				200	-100%	Fails	(±15)	45
EGDN				200	-100%	Fails				200	-100%	Fails	(±15)	

Notes: M = Manually Integrated Signals Differ by More Than 40%  
 D = Operator Disabled Result Signals Differ by More Than 50%  
 O = Over Calibration Range



Inst ID: LC10 Batch ID: 07022009  
Method : Method 8330 Test : SOP SAC-LC-0009  
ICAL Date: See Calibration Report

Date	Time	Operator	Sample ID	File No	Vol or Wt	Extract Vol	Diln	Comments
02-JUL-2009	15:27	fhk	Primer	A-000001.	0 g	0 mL	1	
02-JUL-2009	16:18	fhk	Primer	A-000002.	0 g	0 mL	1	
02-JUL-2009	17:09	fhk	STD_6 09GCSV0054 .4K/.2/.2/.2/	A-000003.	0 g	0 mL	1	
02-JUL-2009	18:00	fhk	LFW791AA 9182192 G9G010000-192	A-000004.	1000 mL	20 mL	1	
02-JUL-2009	18:50	fhk	LFW791AC 9182192 G9G010000-192	A-000005.	1000 mL	20 mL	1	
02-JUL-2009	19:40	fhk	LFQAK1AC 9182192 G9F260326-1	A-000006.	997.78 mL	20 mL	1	
02-JUL-2009	20:31	fhk	LFQAL1AC 9182192 G9F260326-2	A-000007.	991.12 mL	20 mL	1	
02-JUL-2009	21:22	fhk	LFQAM1AC 9182192 G9F260326-3	A-000008.	1009.82 mL	20 mL	1	
02-JUL-2009	22:13	fhk	LFRD61AC 9182192 G9F270189-1	A-000009.	985.56 mL	20 mL	1	
02-JUL-2009	23:03	fhk	LFRD71AC 9182192 G9F270189-2	A-000010.	999.12 mL	20 mL	1	
02-JUL-2009	23:54	fhk	LFRD81AC 9182192 G9F270189-3	A-000011.	991.73 mL	20 mL	1	
03-JUL-2009	00:45	fhk	STD_5 09GCSV0053 .2K/.1/ 1/.1/	A-000012.	0 g	0 mL	1	
03-JUL-2009	01:36	fhk	LFRD91AC 9182192 G9F270189-4	A-000013.	996.78 mL	20 mL	1	
03-JUL-2009	02:27	fhk	LFREA1AC 9182192 G9F270189-5	A-000014.	995.33 mL	20 mL	1	
03-JUL-2009	03:17	fhk	LFV9A1AC 9182192 G9F300242-1	A-000015.	1014.4 mL	20 mL	1	
03-JUL-2009	04:07	fhk	LFV9G1AC 9182192 G9F300242-2	A-000016.	975.28 mL	20 mL	1	
03-JUL-2009	04:58	fhk	LFV9J1AC 9182192 G9F300242-3	A-000017.	1018.21 mL	20 mL	1	
03-JUL-2009	05:49	fhk	LFV9K1AC 9182192 G9F300242-4	A-000018.	1028.11 mL	20 mL	1	
03-JUL-2009	06:39	fhk	LFV9K1AF 9182192 G9F300242-4	A-000019.	1018.28 mL	20 mL	1	
03-JUL-2009	07:30	fhk	LFV9K1AG 9182192 G9F300242-4	A-000020.	1017.3 mL	20 mL	1	
03-JUL-2009	08:20	fhk	LFV9M1AC 9182192 G9F300242-5	A-000021.	1016.72 mL	20 mL	1	
03-JUL-2009	09:11	fhk	STD_6 09GCSV0054 .4K/ 2/.2/.2/	A-000022.	0 g	0 mL	1	
03-JUL-2009	10:01	fhk	09GCSV0234 LEVEL 1	A-000023.	0 g	0 mL	1	
03-JUL-2009	10:52	fhk	09GCSV0235 LEVEL 2	A-000024.	0 g	0 mL	1	
03-JUL-2009	11:42	fhk	09GCSV0236 LEVEL 3	A-000025.	0 g	0 mL	1	
03-JUL-2009	12:33	fhk	09GCSV0237 LEVEL 4	A-000026.	0 g	0 mL	1	
03-JUL-2009	13:23	fhk	09GCSV0238 LEVEL 5	A-000027.	0 g	0 mL	1	
03-JUL-2009	14:14	fhk	09GCSV0239 LEVEL 5	A-000028.	0 g	0 mL	1	
03-JUL-2009	15:04	fhk	09GCSV0240 LEVEL 6	A-000029.	0 g	0 mL	1	
03-JUL-2009	15:55	fhk	09GCSV0241 LEVEL 7	A-000030.	0 g	0 mL	1	
03-JUL-2009	16:45	fhk	09GCSV0242 LEVEL 8	A-000031.	0 g	0 mL	1	
03-JUL-2009	17:36	fhk	09GCSV0243 MRL	A-000032.	0 g	0 mL	1	

Chromatography Summary

Injection Date: 7/2/2009 17:09 Operator: fhk  
 DataFile: LC10.107022009 BVA-000003.D Vial Num: 2  
 Instrument ID: LC10

Method 8330 Target Analyte Results

Sample: **STD\_6 09GCSV0054 .4K/2/2/2/2**

Method File: LC10.107022009 B8330AB.M  
 Start Cal Date: 6/10/2009 15:33 End Cal Date: 6/11/2009 13:47

Matrix: NONE SubList: CAL.sub SpikeList:  
 Samp. Info: STD\_6 09GCSV0054 .4K/2/2/2/2  
 Misc. Info: .6.,.;3;CAL.sub.,0.1

Dilution Factor	Extract Volume	Sample Volume	Sample Weight
1X	0 mL	0 mL	0 g

Signal 1 UV 250-265								Signal 2 UV 358-205								
Compound Name	RT	Response	PPB	Spike Level	%D	Result	Flag	RT	Response	PPB	Spike Level	%D	Result	Flag	Limits(%)	Flag
3,4-Dinitrotoluene	18.20	9636	204.3000<	200	2%	Acceptable		18.20	18445	213.4000	200	7%	Acceptable		(±15)	
HMX	5.29	25432	210.5000<	200	5%	Acceptable					200	-100%	Fails		(±15)	45
RDX	7.85	16295	209.2000<	200	5%	Acceptable					200	-100%	Fails		(±15)	45
Picric ACID	8.99	41811	505.1000	500	1%	Acceptable		8.99	61520	505.6000<	500	1%	Acceptable		(±15)	
1,3,5-Trinitrobenzene	10.20	30367	205.7000<	200	3%	Acceptable					200	-100%	Fails		(±15)	45
1,3-Dinitrobenzene	13.26	29099	208.3000<	200	4%	Acceptable					200	-100%	Fails		(±15)	45
TETRYL	14.32	19114	206.0000<	200	3%	Acceptable					200	-100%	Fails		(±15)	45
Nitrobenzene	15.19	12872	212.1000<	200	6%	Acceptable					200	-100%	Fails		(±15)	45
2,4,6-Trinitrotoluene	16.55	18075	206.1000<	200	3%	Acceptable					200	-100%	Fails		(±15)	45
4-AM-2,6-DNT	17.55	12873	206.6000<	200	3%	Acceptable					200	-100%	Fails		(±15)	45
2-AM-4,6-DNT	18.61	15046	204.5000<	200	2%	Acceptable					200	-100%	Fails		(±15)	45
2,6-Dinitrotoluene	20.34	10970	206.0000<	200	3%	Acceptable					200	-100%	Fails		(±15)	45
2,4-Dinitrotoluene	21.06	17567	205.2000<	200	3%	Acceptable					200	-100%	Fails		(±15)	45
2-Nitrotoluene	24.79	7368	205.0000<	200	3%	Acceptable					200	-100%	Fails		(±15)	45
4-Nitrotoluene	26.61	8982	205.1000<	200	3%	Acceptable					200	-100%	Fails		(±15)	45
3-Nitrotoluene	28.64	8962	207.8000<	200	4%	Acceptable					200	-100%	Fails		(±15)	45
<b>Nitroglycerin</b>				<b>200</b>	<b>-100%</b>	<b>Fails</b>		15.81	12531	206.0000<	200	3%	Acceptable		(±15)	45
<b>PETN</b>				<b>200</b>	<b>-100%</b>	<b>Fails</b>		30.22	6862	214.1000<	200	7%	Acceptable		(±15)	45
3,5-Dinitroaniline	14.01	18502	205.6000<	200	3%	Acceptable					200	-100%	Fails		(±15)	45
<b>EGDN</b>				<b>200</b>	<b>-100%</b>	<b>Fails</b>					200	-100%	Fails		(±15)	

Notes M = Manually Integrated Signals Differ by More Than 40%  
 D = Operator Disabled Result Signals Differ by More Than 50%  
 O = Over Calibration Range

Chromatography Summary

Injection Date: 7/7/2009 0:45 Operator: fhk  
 DataFile: LC10 107022009 BVA-000012.D Vial Num: 3  
 Instrument ID: LC10

Method 8330 Target Analyte Results

Sample: STD\_5 09GCSV0053 .2K/.1/.1/.1

Method File: LC10 107022009 B\8330AB.M  
 Start Cal Date: 6/10/2009 15:33 End Cal Date: 6/11/2009 13:47

Matrix: NONE SubList: CAL sub SpikeList:  
 Samp. Info: STD\_5 09GCSV0053 2K/ 1/ 1/ 1/ 1/ 2  
 Misc. Info: :5; : : 13; CAL sub. :0.1

Dilution Factor	Extract Volume	Sample Volume	Sample Weight
1X	0 mL	0 mL	0 g

Compound Name	RT	Signal 1 UV 250-265						Signal 2 UV 358-205						Limits(%)	Flag
		Response	PPB	Spike Level	%D	Result	Flag	RT	Response	PPB	Spike Level	%D	Result		
3,4-Dinitrotoluene	18.21	4770	101.1000<	100	1%	Acceptable		18.21	9105	105.3000	100	5%	Acceptable	(±15)	
HMX	5.29	12664	104.8000<	100	5%	Acceptable					100	-100%	Fails	(±15)	45
RDX	7.86	8028	103.1000<	100	3%	Acceptable					100	-100%	Fails	(±15)	45
Picric ACID	9.05	16980	205.1000	200	3%	Acceptable	9.05	25011	205.5000<	200	3%	Acceptable	(±15)		
1,3,5-Trinitrobenzene	10.21	15116	102.4000<	100	2%	Acceptable					100	-100%	Fails	(±15)	45
1,3-Dinitrobenzene	13.27	14453	103.5000<	100	4%	Acceptable					100	-100%	Fails	(±15)	45
TETRYL	14.34	9510	102.5000<	100	3%	Acceptable					100	-100%	Fails	(±15)	45
Nitrobenzene	15.20	6322	104.2000<	100	4%	Acceptable					100	-100%	Fails	(±15)	45
2,4,6-Trinitrotoluene	16.57	9018	102.8000<	100	3%	Acceptable					100	-100%	Fails	(±15)	45
4-AM-2,6-DNT	17.56	6403	102.8000<	100	3%	Acceptable					100	-100%	Fails	(±15)	45
2-AM-4,6-DNT	18.63	7508	102.1000<	100	2%	Acceptable					100	-100%	Fails	(±15)	45
2,6-Dinitrotoluene	20.35	5455	102.5000<	100	3%	Acceptable					100	-100%	Fails	(±15)	45
2,4-Dinitrotoluene	21.08	8726	101.9000<	100	2%	Acceptable					100	-100%	Fails	(±15)	45
2-Nitrotoluene	24.79	3635	101.2000<	100	1%	Acceptable					100	-100%	Fails	(±15)	45
4-Nitrotoluene	26.61	4448	101.6000<	100	2%	Acceptable					100	-100%	Fails	(±15)	45
3-Nitrotoluene	28.63	4394	101.9000<	100	2%	Acceptable					100	-100%	Fails	(±15)	45
Nitroglycerin				100	-100%	Fails	15.82	6254	102.8000<	100	3%	Acceptable	(±15)	45	
PETN				100	-100%	Fails	30.25	3304	103.1000<	100	3%	Acceptable	(±15)	45	
3,5-Dinitroaniline	14.02	9179	102.0000<	100	2%	Acceptable					100	-100%	Fails	(±15)	45
EGDN				100	-100%	Fails					100	-100%	Fails	(±15)	

Notes: M = Manually Integrated Signals Differ by More Than 40%  
 D = Operator Disabled Result Signals Differ by More Than 50%  
 O = Over Calibration Range

Chromatography Summary

Injection Date: 7/3/2009 9.11 Operator: shk  
 DataFile: LC10 I07022009 BVA-000022.D Vial Num: 4  
 Instrument ID: LC10

Method 8330 Target Analyte Results

Sample: **STD\_6 09GCSV0054 .4K/2/2/2/2**

Method File: LC10 I07022009.B\8330AB.M  
 Start Cal Date: 6/10/2009 15.33 End Cal Date: 6/11/2009 13.47

Matrix: NONE SubList: CAL.sub SpikeList:  
 Samp. Info: STD\_6 09GCSV0054 4K/2/2/2/2  
 Misc. Info: ,6, , , ,3,CAL.sub, ,0,1

Dilution Factor	Extract Volume	Sample Volume	Sample Weight
1X	0 mL	0 mL	0 g

Signal 1 UV 250-265								Signal 2 UV 358-205								
Compound Name	RT	Response	PPB	Spike Level	%D	Result	Flag	RT	Response	PPB	Spike Level	%D	Result	Flag	Limits(%)	Flag
3,4-Dinitrotoluene	18.24	9589	203.3000<	200	2%	Acceptable		18.23	18290	211.6000	200	6%	Acceptable		(±15)	
HMX	5.30	25440	210.6000<	200	5%	Acceptable					200	-100%	Fails		(±15)	45
RDX	7.87	16268	208.8000<	200	4%	Acceptable					200	-100%	Fails		(±15)	45
Picric ACID	9.04	42655	515.3000	500	3%	Acceptable		9.04	62723	515.5000<	500	3%	Acceptable		(±15)	
1,3,5-Trinitrobenzene	10.22	30282	205.1000<	200	3%	Acceptable					200	-100%	Fails		(±15)	45
1,3-Dinitrobenzene	13.29	28991	207.5000<	200	4%	Acceptable					200	-100%	Fails		(±15)	45
TETRYL	14.35	19191	206.8000<	200	3%	Acceptable					200	-100%	Fails		(±15)	45
Nitrobenzene	15.21	12738	209.9000<	200	5%	Acceptable					200	-100%	Fails		(±15)	45
2,4,6-Trinitrotoluene	16.58	18022	205.5000<	200	3%	Acceptable					200	-100%	Fails		(±15)	45
4-AM-2,6-DNT	17.58	12828	205.9000<	200	3%	Acceptable					200	-100%	Fails		(±15)	45
2-AM-4,6-DNT	18.65	15059	204.7000<	200	2%	Acceptable					200	-100%	Fails		(±15)	45
2,6-Dinitrotoluene	20.38	10882	204.4000<	200	2%	Acceptable					200	-100%	Fails		(±15)	45
2,4-Dinitrotoluene	21.10	17483	204.2000<	200	2%	Acceptable					200	-100%	Fails		(±15)	45
2-Nitrotoluene	24.82	7294	203.0000<	200	2%	Acceptable					200	-100%	Fails		(±15)	45
4-Nitrotoluene	26.64	8676	202.6000<	200	1%	Acceptable					200	-100%	Fails		(±15)	45
3-Nitrotoluene	28.67	8836	204.9000<	200	2%	Acceptable					200	-100%	Fails		(±15)	45
Nitroglycerin				200	-100%	Fails		15.84	12481	205.2000<	200	3%	Acceptable		(±15)	45
PETN				200	-100%	Fails		30.27	6681	208.4000<	200	4%	Acceptable		(±15)	45
3,5-Dinitroaniline	14.03	18458	205.1000<	200	3%	Acceptable					200	-100%	Fails		(±15)	45
EGDN				200	-100%	Fails					200	-100%	Fails		(±15)	

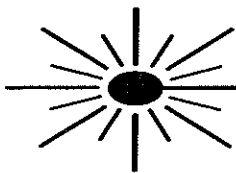
Notes: M = Manually Integrated Signals Differ by More Than 40%  
 D = Operator Disabled Result Signals Differ by More Than 50%  
 O = Over Calibration Range

**Data Validation Report**

**Michael Baker Jr., Inc**

**Camp Bonneville**

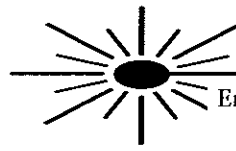
**SDG#: PSF0839**



**ECT.CON INC.**

Environmental and Computer  
Technology Consultants

3531 Fox Chase Drive  
Imperial, PA 15126  
(724) 695-8042  
FAX (724) 695-2698  
e-mail: [ectconinc@comcast.net](mailto:ectconinc@comcast.net)



## Data Validation Report

SDG#	PSF0839
Validation Report Date	August 17, 2009
Validation Guidance	USEPA CLP National Functional Guidelines for Data Review
Client Name	Michael Baker Jr., Inc.
Project Name	Camp Bonneville
Laboratory	TestAmerica
Analytical Parameters	Explosives
Analytical Method	SW-846 8330

### Samples/Matrix:

Date Sampled	Location	Portland ID	Sacramento ID	Explosives	Matrix
06/24/09	23LCMW03DW	PSF0839-01	LFRD6	X	Aqueous
06/25/09	23LCMW03SW	PSF0839-02	LFRD7	X	Aqueous
06/25/09	23LCMW02DW	PSF0839-03	LFRD8	X	Aqueous
06/25/09	23LCMW02SW	PSF0839-04	LFRD9	X	Aqueous
06/25/09	23LCMW01DW	PSF0839-05	LFREA	X	Aqueous

Analytical data in this report were screened to determine analytical limitations of the data based on specific quality control criteria. This screening assumes analytical results are correct as reported and merely provides an interpretation of the reported quality control results. Laboratory calculations have been verified as part of this validation. Specific findings on analytical limitations are presented in this report. Annotated Form 1s or spreadsheets for samples reviewed are included after the Data Assessment Findings. Form 1s for the MS/MSD samples and spreadsheets are not annotated.

### SUMMARY

The sample set for the Camp Bonneville site consists of five aqueous field samples. These samples were analyzed for explosives as provided in the Table 1. The findings presented in this review of the analytical data assume that the information presented by the analytical laboratory is correct.

The explosives findings are based upon the assessment of the following:

- \* ● Data Completeness
- \* ● Holding Times
- \* ● Calibration (Initial and Continuing)
- Blanks
- \* ● System Monitoring Compounds (Surrogate Spikes)
- \* ● Matrix Spike/Matrix Spike Duplicates
- \* ● Laboratory Control Standard Results
- \* ● Target Compound Identification
- Compound Quantification and Reported Contract Quantitation Limits

- \* • System Performance
- \* Criteria were met for this evaluation item.

This evaluation was conducted in accordance with USEPA CLP National Functional Guidelines for Organic Data Review and the analytical method. Findings from this evaluation should be considered when using the analytical data. This report presents a summary of the data qualifications based on the review of the aforementioned evaluation criteria. This is followed by annotated Form 1s/ spreadsheets. Finally, the worksheets used to perform the evaluation are provided.

## **FINDINGS**

### **1. Blanks**

The laboratory method blank exhibited contamination for the following parameter:

Blank	Compound	Maximum Concentration (ppb)	Action Limit (ppm)	Action
LFW791	1,3,5-Trinitrobenzene	0.038	0.19	U sample results < RL

RL - reporting limit

### **2. Compound Quantitation**

In sample 23LCMW02DW, the positive result below the reporting limit for 2-nitrotoluene was not confirmed. Therefore, the form one was amended to show the result as nondetected "U" at the reporting limit.

## **NOTES**

### **Matrix Spike/Matrix Spike Duplicate Results**

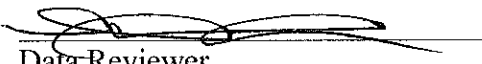
A MS/MSD was not analyzed with this SDG. A laboratory control sample was analyzed. Data were not qualified on this basis.

### **Laboratory Control Sample Results**

Recovery of 12 out of 17 compounds exceeded the laboratory's upper quality control limits. The affected compounds were not detected in the associated samples. Data were not qualified on this basis.

### **Field Duplicate Results**

No field duplicates were included with this SDG. Data were not qualified on this basis.

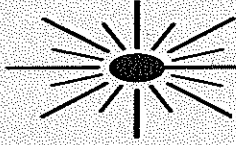
  
Data Reviewer

8/17/09  
Date



### **Glossary of Data Qualifiers**

U	Not Detected.	The associated number indicates approximate sample concentration necessary to be detected.
UJ	Not Detected.	Quantitation limit may be inaccurate or imprecise.
J	Analyte Present.	Reported value may not be accurate or precise.
N	Consider Present.	Tentative identification. Special methods may be needed to confirm its presence or absence in future sampling efforts.
R	Unusable Result.	Analyte may or may not be present in the sample.
UR	Unusable Result.	Analyte may or may not be present in the sample.



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*Annotated Form 1's*  
*(Spreadsheet)*

TestAmerica Portland

Client Sample ID: PSF0839-01

HPLC

23LCMW030W

Lot-Sample #...: G9F270189-001    Work Order #...: LFRD61AC    Matrix.....: WATER  
 Date Sampled...: 06/24/09    Date Received...: 06/27/09  
 Prep Date.....: 07/01/09    Analysis Date...: 07/02/09  
 Prep Batch #...: 9182192  
 Dilution Factor: 1.01    Method.....: SW846 8330

PARAMETER	RESULT	REPORTING		
		LIMIT	UNITS	MDL
Nitroglycerin	ND	0.66	ug/L	0.15
PETN	ND	0.66	ug/L	0.23
Picric Acid	ND	1.0	ug/L	0.13
2-Amino-4,6-dinitrotoluene	ND	0.20	ug/L	0.10
4-Amino-2,6-dinitrotoluene	ND	0.10	ug/L	0.022
1,3-Dinitrobenzene	ND	0.10	ug/L	0.050
2,4-Dinitrotoluene	ND	0.10	ug/L	0.050
2,6-Dinitrotoluene	ND	0.10	ug/L	0.050
HMX	ND	0.10	ug/L	0.027
Nitrobenzene	ND	0.10	ug/L	0.050
2-Nitrotoluene	ND	0.40	ug/L	0.073
3-Nitrotoluene	ND	0.40	ug/L	0.063
4-Nitrotoluene	ND	0.50	ug/L	0.073
RDX	ND	0.10	ug/L	0.066
Tetryl	ND	0.10	ug/L	0.050
1,3,5-Trinitrobenzene	<del>0.035-J,B</del>	0.10	ug/L	0.031
2,4,6-Trinitrotoluene	ND	0.10	ug/L	0.024
	PERCENT	RECOVERY		
SURROGATE	RECOVERY	LIMITS		
3,4-Dinitrotoluene	105	(79 - 111)		

NOTE(S):

J Estimated result. Result is less than RL.

B Method blank contamination. The associated method blank contains the target analyte at a reportable level.

*WJ*  
*8/10/09*

TestAmerica Portland

Client Sample ID: PSF0839-02

HPLC

23CMW 035W

Lot-Sample #...: G9F270189-002 Work Order #...: LFRD71AC Matrix.....: WATER  
 Date Sampled...: 06/25/09 Date Received...: 06/27/09  
 Prep Date.....: 07/01/09 Analysis Date...: 07/02/09  
 Prep Batch #...: 9182192  
 Dilution Factor: 1 Method.....: SW846 8330

PARAMETER	RESULT	REPORTING LIMIT	UNITS	MDL
Nitroglycerin	ND	0.65	ug/L	0.15
PETN	ND	0.65	ug/L	0.23
Picric Acid	ND	1.0	ug/L	0.12
2-Amino-4,6-dinitrotoluene	ND	0.20	ug/L	0.10
4-Amino-2,6-dinitrotoluene	ND	0.10	ug/L	0.022
1,3-Dinitrobenzene	ND	0.10	ug/L	0.050
2,4-Dinitrotoluene	ND	0.10	ug/L	0.050
2,6-Dinitrotoluene	ND	0.10	ug/L	0.050
HMX	ND	0.10	ug/L	0.027
Nitrobenzene	ND	0.10	ug/L	0.050
2-Nitrotoluene	ND	0.40	ug/L	0.072
3-Nitrotoluene	ND	0.40	ug/L	0.062
4-Nitrotoluene	ND	0.50	ug/L	0.072
RDX	ND	0.10	ug/L	0.065
Tetryl	ND	0.10	ug/L	0.050
1,3,5-Trinitrobenzene	<del>0.037 J, B</del>	0.10 U	ug/L	0.031
2,4,6-Trinitrotoluene	ND	0.10	ug/L	0.024

SURROGATE	PERCENT RECOVERY	RECOVERY LIMITS
3,4-Dinitrotoluene	106	(79 - 111)

NOTE(S):

- J Estimated result Result is less than RL.
- B Method blank contamination. The associated method blank contains the target analyte at a reportable level

WKS  
8/6/09

TestAmerica Portland

Client Sample ID: PSF0839-03

HPLC *23LCMW02DW*

Lot-Sample #...: G9F270189-003    Work Order #...: LFRD81AC    Matrix.....: WATER  
 Date Sampled...: 06/25/09    Date Received...: 06/27/09  
 Prep Date.....: 07/01/09    Analysis Date...: 07/02/09  
 Prep Batch #...: 9182192  
 Dilution Factor: 1    Method.....: SW846 8330

PARAMETER	RESULT	REPORTING		
		LIMIT	UNITS	MDL
Nitroglycerin	ND	0.65	ug/L	0.15
PETN	ND	0.65	ug/L	0.23
Picric Acid	ND	1.0	ug/L	0.12
2-Amino-4,6-dinitrotoluene	ND	0.20	ug/L	0.10
4-Amino-2,6-dinitrotoluene	ND	0.10	ug/L	0.022
1,3-Dinitrobenzene	ND	0.10	ug/L	0.050
2,4-Dinitrotoluene	ND	0.10	ug/L	0.050
2,6-Dinitrotoluene	ND	0.10	ug/L	0.050
HMX	ND	0.10	ug/L	0.027
Nitrobenzene	ND	0.10	ug/L	0.050
2-Nitrotoluene	<del>0.20</del> <i>JJ</i>	0.40 <i>U</i>	ug/L	0.072
3-Nitrotoluene	ND	0.40	ug/L	0.062
4-Nitrotoluene	ND	0.50	ug/L	0.072
RDX	ND	0.10	ug/L	0.065
Tetryl	ND	0.10	ug/L	0.050
1,3,5-Trinitrobenzene	<del>0.034</del> <i>J,B</i>	0.10 <i>U</i>	ug/L	0.031
2,4,6-Trinitrotoluene	ND	0.10	ug/L	0.024
	PERCENT	RECOVERY		
<u>SURROGATE</u>	<u>RECOVERY</u>	<u>LIMITS</u>		
3,4-Dinitrotoluene	104	(79 - 111)		

NOTE(S):

- J Estimated result. Result is less than RL.
- B Method blank contamination. The associated method blank contains the target analyte at a reportable level.

*WJ  
8/10/09*

TestAmerica Portland

Client Sample ID: PSF0839-04

HPLC

23LCMW02SW

Lot-Sample #...: G9F270189-004 Work Order #...: LFRD91AC Matrix.....: WATER  
 Date Sampled...: 06/25/09 Date Received...: 06/27/09  
 Prep Date.....: 07/01/09 Analysis Date...: 07/03/09  
 Prep Batch #...: 9182192  
 Dilution Factor: 1 Method.....: SW846 8330

PARAMETER	RESULT	REPORTING		
		LIMIT	UNITS	MDL
Nitroglycerin	ND	0.65	ug/L	0.15
PETN	ND	0.65	ug/L	0.23
Picric Acid	ND	1.0	ug/L	0.12
2-Amino-4,6-dinitrotoluene	ND	0.20	ug/L	0.10
4-Amino-2,6-dinitrotoluene	ND	0.10	ug/L	0.022
1,3-Dinitrobenzene	ND	0.10	ug/L	0.050
2,4-Dinitrotoluene	ND	0.10	ug/L	0.050
2,6-Dinitrotoluene	ND	0.10	ug/L	0.050
HMX	ND	0.10	ug/L	0.027
Nitrobenzene	ND	0.10	ug/L	0.050
2-Nitrotoluene	ND	0.40	ug/L	0.072
3-Nitrotoluene	ND	0.40	ug/L	0.062
4-Nitrotoluene	ND	0.50	ug/L	0.072
RDX	ND	0.10	ug/L	0.065
Tetryl	ND	0.10	ug/L	0.050
1,3,5-Trinitrobenzene	<del>0.037</del> J,B	0.10 U	ug/L	0.031
2,4,6-Trinitrotoluene	ND	0.10	ug/L	0.024
	PERCENT	RECOVERY		
SURROGATE	RECOVERY	LIMITS		
3,4-Dinitrotoluene	105	(79 - 111)		

NOTE(S):

- J Estimated result. Result is less than RL.
- B Method blank contamination. The associated method blank contains the target analyte at a reportable level

WJG  
8/16/09

TestAmerica Portland

Client Sample ID: PSF0839-05

HPLC *23LCMWO1DW*

Lot-Sample #...: G9F270189-005    Work Order #...: LFREA1AC    Matrix.....: WATER  
 Date Sampled...: 06/25/09    Date Received...: 06/27/09  
 Prep Date.....: 07/01/09    Analysis Date...: 07/03/09  
 Prep Batch #...: 9182192  
 Dilution Factor: 1.01    Method.....: SW846 8330

PARAMETER	RESULT	REPORTING LIMIT	UNITS	MDL
Nitroglycerin	ND	0.66	ug/L	0.15
PETN	ND	0.66	ug/L	0.23
Picric Acid	ND	1.0	ug/L	0.13
2-Amino-4,6-dinitrotoluene	ND	0.20	ug/L	0.10
4-Amino-2,6-dinitrotoluene	ND	0.10	ug/L	0.022
1,3-Dinitrobenzene	ND	0.10	ug/L	0.050
2,4-Dinitrotoluene	ND	0.10	ug/L	0.050
2,6-Dinitrotoluene	ND	0.10	ug/L	0.050
HMX	ND	0.10	ug/L	0.027
Nitrobenzene	ND	0.10	ug/L	0.050
2-Nitrotoluene	ND	0.40	ug/L	0.073
3-Nitrotoluene	ND	0.40	ug/L	0.063
4-Nitrotoluene	ND	0.50	ug/L	0.073
RDX	ND	0.10	ug/L	0.066
Tetryl	ND	0.10	ug/L	0.050
1,3,5-Trinitrobenzene	<del>0.038</del> <i>JTB</i>	0.10 <i>U</i>	ug/L	0.031
2,4,6-Trinitrotoluene	ND	0.10	ug/L	0.024
	PERCENT	RECOVERY		
SURROGATE	RECOVERY	LIMITS		
3,4-Dinitrotoluene	105	(79 - 111)		

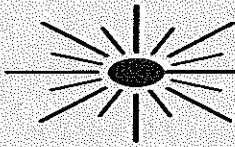
NOTE(S) :

J Estimated result. Result is less than RL.

B Method blank contamination. The associated method blank contains the target analyte at a reportable level.

*LUK  
8/6/09*





ECT.CON INC.

Environmental and Computer  
Technology Consultants

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## *Support Documentation*

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*Laboratory Case Narrative*



**TestAmerica**

THE LEADER IN ENVIRONMENTAL TESTING

## **DATA DELIVERABLES PACKAGE**

Michael Baker Jr., Inc.  
James D. Peyton  
5261 Fountain Drive, Suite A  
Crown Point, IN 46307

Client Project: Camp Bonneville Groundwater  
Client Project Number: 110002 Task 6200

TA Work Order#: PSF0839  
TA Project Manager: Estella Rieben

The total number of pages contained in this data package is:

154

July 31, 2009

TestAmerica - Portland  
9405 S. W. Nimbus Avenue  
Beaverton, Oregon 97008  
(503) 906-9200  
(503) 906-9210



**Michael Baker Jr., Inc.**

5261 Fountain Drive, Suite A  
Crown Point, IN 46307

Project Name: **Camp Bonneville Groundwater**  
Project Number: 110002 Task 6200  
Project Manager: James D. Peyton

Report Created:  
07/21/09 08:43

## ANALYTICAL REPORT FOR SAMPLES

Sample ID	Laboratory ID	Matrix	Date Sampled	Date Received
23LCMW03DW	PSF0839-01	Water	06/24/09 17:45	06/25/09 17:45
23LCMW03SW	PSF0839-02	Water	06/25/09 09:30	06/25/09 17:45
23LCMW02DW	PSF0839-03	Water	06/25/09 11:30	06/25/09 17:45
23LCMW02SW	PSF0839-04	Water	06/25/09 13:00	06/25/09 17:45
23LCMW01DW	PSF0839-05	Water	06/25/09 14:45	06/25/09 17:45
TB261	PSF0839-06	Water	06/25/09 00:00	06/25/09 17:45

TestAmerica Portland



Estella Rieben, Project Manager

*The results in this report apply to the samples analyzed in accordance with the chain of custody document. This analytical report shall not be reproduced except in full, without the written approval of the laboratory.*



# TestAmerica

THE LEADER IN ENVIRONMENTAL TESTING

July 28, 2009

**TestAmerica Project Number: G9F270189**  
PO/Contract: PSF0839

Estella Rieben  
TestAmerica Portland  
Nimbus Corporate Center  
9405 SW Nimbus Ave  
Beaverton, OR 97008

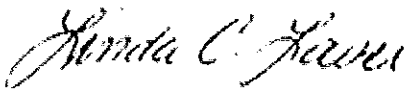
Dear Ms. Rieben,

This report contains the analytical results for the samples received under chain of custody by TestAmerica on June 27, 2009. These samples are associated with your Michael Baker Jr., Inc. project.

The test results in this report meet all NELAC requirements for parameters that accreditation is required or available. Any exceptions to NELAC requirements are noted in the case narrative. The case narrative is an integral part of this report.

If you have any questions, please feel free to call me at (916) 374-4362.

Sincerely,



Linda C. Laver  
Project Manager

## Case Narrative

### TestAmerica West Sacramento Project Number G9F270189

#### General Comments

One of the two bottles for sample PSF0839-01 was received broken. One of the two bottles for sample PSF0839-02 was received with a broken cap. The cap was replaced and this bottle was marked for backup use only.

#### **WATER, 8330, Nitroaromatics & Nitramines**

Samples: 1, 2, 3, 4, 5

The laboratory control sample (LCS) associated with this extraction batch has percent recoveries for most analytes above the established control limits indicating a potential high bias in the data. These samples do not have detected concentrations above the reporting limit for these analytes and there is no adverse impact upon the data.

The matrix spike/matrix spike duplicate (MS/MSD) associated with this extraction batch also has recoveries and/or precision is outside the established control limits for many analytes. This anomaly is most likely matrix related.

There are no other anomalies associated with this project.

## Sample Summary

### TestAmerica West Sacramento Project Number G9F270189

<u>WO#</u>	<u>Sample #</u>	<u>Client Sample ID</u>	<u>Sampling Date</u>	<u>Received Date</u>
LFRD6	1	PSF0839-01	6/24/2009 05:45 PM	6/27/2009 10:00 AM
LFRD7	2	PSF0839-02	6/25/2009 09:30 AM	6/27/2009 10:00 AM
LFRD8	3	PSF0839-03	6/25/2009 11:30 AM	6/27/2009 10:00 AM
LFRD9	4	PSF0839-04	6/25/2009 01:00 PM	6/27/2009 10:00 AM
LFREA	5	PSF0839-05	6/25/2009 02:45 PM	6/27/2009 10:00 AM

#### Notes(s):

- The analytical results of the samples listed above are presented on the following pages.
- All calculations are performed before rounding to avoid round-off errors in calculated results.
- Results noted as "ND" were not detected at or above the stated limit.
- This report must not be reproduced, except in full, without the written approval of the laboratory.
- Results for the following parameters are never reported on a dry weight basis: color, corrosivity, density, flashpoint, ignitability, layers, odor, paint filter test, pH, porosity, pressure, reactivity, redox potential, specific gravity, spot tests, solids, solubility, temperature, viscosity, and weight.



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# *Chain of Custody*

# TestAmerica

THE LEADER IN ENVIRONMENTAL TESTING

11720 North Creek Pkwy N, Suite 400, Bothell, WA 98011-8244  
 1922 E. First Ave., Spokane, WA 99206-5400  
 2405 SW Nimbus Ave, Beaverton, OR 97008-7147  
 100 W. Satehill Ave, #100, Redwood City, CA 94061-1100

254-420-0290 FAX 420-0210  
 509-477-0280 FAX 477-0200  
 503-970-0280 FAX 970-0200  
 415-992-0280 FAX 992-0200

## CHAIN OF CUSTODY REPORT

Work Order #: **PSF 0839**

CLIENT: Michael Baker Jr., Inc  
 REPORT TO: Margaret James  
 ADDRESS: 5761 Forest Hill Drive S.W.  
 Clatsop County, OR 97130  
 PHONE: 503-736-0263 FAX: 503-736-0263

INVOICE TO: Baker

PROJECT NAME: Camp Beineville  
 PROJECT NUMBER: 110003 T 6000

SAMPLED BY: Margaret James

TURNAROUND REQUEST  
 in Business Days  
 Organic & Inorganic Analysis  
 Repetition: 1 2 3 4 5  
 Method: 1 2 3 4 5

OTHER: Specify: \_\_\_\_\_

\* Turnaround Request is their standard may vary by lab charges.

CLIENT SAMPLE IDENTIFICATION	SAMPLING DATE/TIME	HCL	NA	NA	NA	NA	NA	NA	NA	MATRIX (W, S, O)	# OF CONT.	LOCATION/ COMMENTS	TA	WO ID
23LCMW03DW	6/24/09-1745	X	X	X	X	X	X	X	X	W	6	NoTICS		
23LCMW03SW	6/25/09-0930	X	X	X	X	X	X	X	X	W	6	NoTICS		
23LCMW02DW	6/25/09-1130	X	X	X	X	X	X	X	X	W	6	NoTICS		
23LCMW02SW	6/25/09-1300	X	X	X	X	X	X	X	X	W	6	NoTICS		
23LCMW01DW	6/25/09-1445	X	X	X	X	X	X	X	X	W	6	NoTICS		
T0261	6/25/09	X								W	1	NoTICS		
7														
8														
9														
10														

RECEIVED BY: Baker  
 PRINT NAME: Baker  
 DATE: 6/25/09  
 TIME: 1600

RECEIVED BY: Margaret James  
 PRINT NAME: Margaret James  
 DATE: 6/25/09  
 TIME: 1745

ADDITIONAL REMARKS: TAP

DATE: 6/25/09  
 TIME: 1612  
 DATE: 6/25/09  
 TIME: 1745

TEMP: 34

PAGE: 34

TestAmerica Portland  
**Sample Receiving Checklist**

Work Order #: PSF0839 Date/Time Received: 4/25/09 1745  
 Client Name and Project: Michael Baker Camp Bonnaville

Time Zone:  
 EDT/EST     CDT/CST     MDT/MST     PDT/PST     AK     OTHER

**Unpacking Checks:**

Cooler #(s): 1  
 Temperatures: 3.4  
 Digi #1  Digi #2  IR Gun Raytek  
 Plastic  Glass

**Temperature out of Range:**

Not enough or No Ice  
 Ice Melted  
 W/in 4 Hrs of collection  
 Other: \_\_\_\_\_

- N/A Yes No Initials: BIE
1. If ESI client, were temp blanks received? If no, document on NOD.
  2. Cooler Seals intact? (N/A if hand delivered) if no, document on NOD.
  3. Chain of Custody present? If no, document on NOD.
  4. Bottles received intact? If no, document on NOD.
  5. Sample is not multiphasic? If no, document on NOD.
  6. Proper Container and preservatives used? If no, document on NOD.
  7. pH of all samples checked and meet requirements? If no, document on NOD.
  8. Cyanide samples checked for sulfides and meet requirements? If no, notify PM.
  9. HF Dilution required?
  10. Sufficient volume provided for all analysis? If no, document on NOD and consult PM before proceeding.
  11. Did chain of custody agree with samples received? If no, document on NOD.
  12. Is the "Sampled by" section of the COC completed?
  13. Were VOA/Oil Syringe samples without headspace?
  14. Were VOA vials preserved?  PCl  Sodium Thiosulfate  Ascorbic Acid
  15. Did samples require preservation with sodium thiosulfate?
  16. If yes to #14, was the residual chlorine test negative? If no, document on NOD.
  17. Are dissolved/field filtered metals bottles sediment-free? If no, document on NOD.
  18. Is sufficient volume provided for client requested MS/MSD or matrix duplicates? If no, document on NOD and contact PM before proceeding.
  19. Are analyses with short holding times received in hold?
  20. Was Standard Turn Around (TAT) requested?
  21. Receipt date(s) < 48 hours past the collection date(s)? If no, notify PM.

TestAmerica Portland  
Sample Receiving Checklist

Work Order #: PSF0839

Login Checks:

Initials: PS

N/A Yes No

22. Sufficient volume provided for all analysis? If no, document on NOD & contact PM.
23. Sufficient volume provided for client requested MS/MSD or matrix duplicates? If no, document on NOD and contact PM.
24. Did the chain of custody include "received by" and "relinquished by" signatures, dates and times?
25. Were special log in instructions read and followed?
26. Were tests logged checked against the COC?
27. Were rush notices printed and delivered?
28. Were short hold notices printed and delivered?
29. Were subcontract COCs printed?
30. Was HF dilution logged?

Labeling and Storage Checks:

Initials: MM

N/A Yes No

31. Were the subcontracted samples/containers put in Sx fridge?
32. Were sample bottles and COC double checked for dissolved/filtered metals?
33. Did the sample ID, Date, and Time from label match what was logged?
34. Were Foreign sample stickers affixed to each container and containers stored in foreign fridge?
35. Were HF stickers affixed to each container, and containers stored in Sx fridge?
36. Was an NOD for created for noted discrepancies and placed in folder?

Document any problems or discrepancies and the actions taken to resolve them on a Notice of Discrepancy form (NOD).

CLIENT TAL - Portland PM LL LOG# 59241  
 LOT# (QUANTIMS ID) GAF270189 QUOTE# 83365 LOCATION W21C

DATE RECEIVED 6-27-09 TIME RECEIVED 1000 Initials AK Date 6-27-09

DELIVERED BY  FEDEX  CA OVERNIGHT  CLIENT  
 AIRBORNE  GOLDENSTATE  DHL  
 UPS  BAX GLOBAL  GO-GETTERS  
 TAL COURIER  VALLEY LOGISTICS  MORGAN HILL COURIER  
 OTHER

CUSTODY SEAL STATUS  INTACT  BROKEN  N/A

CUSTODY SEAL #(S) 204265

SHIPPING CONTAINER(S)  TAL  CLIENT  N/A

TEMPERTURE RECORD (IN °C) 4  5  OTHER

COC #(S) N/A

TEMPERATURE BLANK Observed: 0 Corrected: \_\_\_\_\_

SAMPLE TEMPERATURE  
 Observed: 0 Average: 0 Corrected Average: 0

COLLECTOR'S NAME  Verified from COC  Not on COC

pH MEASURED  YES  ANOMALY  N/A

LABELED BY.....

LABELS CHECKED BY.....

PEER REVIEW  N/A

SHORT HOLD TEST NOTIFICATION SAMPLE RECEIVING  
 WETCHEM  N/A  
 VOA-ENCORES  N/A

METALS NOTIFIED OF FILTER/PRESERVE VIA VERBAL & EMAIL  N/A

COMPLETE SHIPMENT RECEIVED IN GOOD CONDITION WITH APPROPRIATE TEMPERATURES, CONTAINERS, PRESERVATIVES  N/A

CLOUSEAU  TEMPERATURE EXCEEDED (2 °C - 6 °C)\*  N/A

WET ICE  BLUE ICE  GEL PACK  NO COOLING AGENTS USED  PM NOTIFIED

Notes: PSF 0839-01 - rec'd 1 of 2 ASB broken  
" " -02 - rec'd 1 bottle cap broken & replaced. idrot  
no notice on bottle.

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# *Worksheets*

**HOLDING TIMES**

SAMPLE DATE	SAMPLE ID	Portland LAB ID	Sacramento Lab ID	Matrix	Prep	Anal
06/24/09	23LCMW03DW	PSF0839-01	LFRD6	Aq	07/01/09	7/2/09 1X
06/25/09	23LCMW03SW	PSF0839-02	LFRD7	Aq	07/01/09	7/2/09 1X
06/25/09	23LCMW02DW	PSF0839-03	LFRD8	Aq	07/01/09	7/2/09 1X
06/25/09	23LCMW02SW	PSF0839-04	LFRD9	Aq	07/01/09	7/3/09 1X
06/25/09	23LCMW01DW	PSF0839-05	LFREA	Aq	07/01/09	7/3/09 1X

7 days from collection to extraction  
 Sample Date 6/24/2009 Extract By 7/1/2009  
 40 days from extraction to analysis  
 Sample Date 7/1/2009 Extract By 8/10/2009

**TARGET COMPOUNDS AND QUANTITATION LIMITS**

Was a target compound list provided by the client? No  
 Did Sample Form 1s match the target compound list NA  
 Were required quantitation limits provided by the client? No  
 Did all compounds meet the required quantitation limits? NA

**SYSTEM MONITORING COMPOUNDS**

SAMPLE	SURROGATE	COLUMN 1	DF	ACTION
All IN				

Were surrogate RTs within windows established by the ICAL? Yes  
 Were there any transcription errors between the raw data and Form 2? No  
 Were laboratory acceptance limits used as the basis for validation? 30-150  
 Did the laboratory provide CLP Form II or equivalent? Yes

02DW

SURR	AMOUNT FOUND	AMOUNT SPIKED	% R	F1
34-DNT A			#DIV/0!	104
34-DNT C			#DIV/0!	

**MS/MSD RECOVERY and RELATIVE PERCENT DIFFERENCE**

Non-Client

**LABORATORY CONTROL SAMPLES**

SAMPLE	COMPOUND	LCS %R	LCSD %R	RPD	ACTION
LFW791	12 of 17 HI		NA	NA	none, ND

Were laboratory acceptance limits used as the basis for validation? Yes  
 Did the laboratory provide CLP Form III or equivalent? Yes  
 Were chromatograms and quan reports present for all LCS/LCDs? Yes

LFW791

RDX

12 of 17 HI	AMOUNT FOUND	AMOUNT SPIKED	% R	FORM 3
LCS	127	1	127.0	127

%R = (Amount Found/Amount Spiked)\*100

**BLANKS**

BLANK	COMPOUND	RESULT	5X OR 10X	ACTION LEVEL	ACTION	
LFW791	135-TNBENZ	0.038	5	0.19	U at RL	ALL

NOTE Equipment/field blanks are not qualified on the basis of laboratory method blank contamination or contamination in other field quality control blanks.

**CALIBRATION**

**Initial**

Are chromatograms and quan reports present for all ICAL standards? Yes  
 Are CLP For VIs or equivalent present and complete? Yes  
 Any transcription or calculation errors? No  
 What linearity criteria was used? 20% RSD or 0.99  
 Were RT window documented? Yes

**Continuing**

Are chromatograms and quan reports present for all CCV standards? Yes  
 Are CLP Form VIs or equivalent present and complete? Yes  
 Any transcription or calculation errors? No  
 What %D criteria was used? 25% D  
 Have all samples and standards been listed on an Analytical Sequence? Yes  
 Was a proper analytical sequence followed? Yes

**INITIAL CALIBRATION - A**

Date	6/10/2009	Rptd Avg CF	77.89163	Rptd %RSD	5.156
Instrument	A	Cald Avg CF	77.892	Cald %RSD	5.156
Compound	RDX				
Reported CF	80.13	CF1	81.80	STD DEV	4
Calculated CF	80.42	CF2	75.90		
		CF3	77.15		
		CF4	80.72		
Response	16084	CF5	78.45		
Conc	200	CF6	80.13		
		CF7	69.168	%RSD = (Std Dev/Avg CF)*100	
	CF = (H/C)	CF8	79.815		

**CONTINUING CALIBRATION A**

Date	7/2/2009		
Time	1709		
Instrument	A		
Compound	HMX		
Reported Conc	104.5	Rptd %D	5
Calculated Conc	210.182	Cald %D	-5.09
Response	25432	ICAL CF	200
CF	121	CCV CF	210.2
	CF = (H/C)	%Difference = ((ICAL - CCV)/ICAL)*100	

**FIELD DUPLICATES**

COMPOUND	SAMPLE	QUALIFIER	DUPLICATE	QUALIFIER	RPD
PCA					#DIV/0!

NOTES Samples are not qualified on this basis.

**COMPOUND IDENTIFICATION AND QUANTITATION**

Has a F10 been completed for every sample containing positive results? Yes  
 Was RT data presented on the form? NA  
 Are RTs within the established windows? Yes  
 Any transcription or calculation errors? No  
 Any false positives, negative peaks, shouldering, etc.? No  
 Was GCMS confirmation needed for results > 10 µg/ml? NA  
 Were percent differences or relative percent differences calculated? NA  
 Are percent differences/RPDs greater than 25%? NA  
 Are there any transcription errors? No  
 Are Form Is present for all field and quality control samples? No  
 Are chromatograms and quan reports present for all samples? Yes  
 Are RLs adjusted to reflect sample dilutions, percent solids, etc.? Yes  
 For soils, any percent solids <50%? No  
 For soils, any percent solids <10%? No

23L4MW01AW RDX 0.13 µg/L

	A	C
response	524	269
cf	77.89163	42.55812
final vol ml	20	20
initial L	1.02157	1.02157
df	1	1
calculated	0.1317	0.1237

**PERCENT DIFFERENCE BETWEEN COLUMNS**

SAMPLE	COMPOUND	A	C	RPD	QUALIFIER
23LCMW02DW	2-ntoluene	0.1953	not confirmed	#VALUE!	U at RL



Non-Client

MATRIX SPIKE SAMPLE EVALUATION REPORT

HPLC

Client Lot #...: G9F270189      Work Order #...: LFV9K1AF-MS      Matrix.....: WATER  
MS Lot-Sample #: G9F300242-004      LFV9K1AG-MSD  
Date Sampled...: 06/26/09      Date Received...: 06/30/09  
Prep Date.....: 07/01/09      Analysis Date...: 07/03/09  
Prep Batch #...: 9182192  
Dilution Factor: 0.98

PARAMETER	PERCENT RECOVERY	RECOVERY LIMITS	RPD	RPD LIMITS	METHOD
Nitroglycerin	106	(85 - 115)			SW846 8330
	108	(85 - 115)	2.0	(0-15)	SW846 8330
PEIN	110	(84 - 117)			SW846 8330
	111	(84 - 117)	0.90	(0-15)	SW846 8330
Picric Acid	70	(21 - 118)			SW846 8330
	99 p	(21 - 118)	35	(0-20)	SW846 8330
2-Amino-4,6-dinitrotoluene	125 a	(77 - 123)			SW846 8330
	129 a	(77 - 123)	3.4	(0-27)	SW846 8330
4-Amino-2,6-dinitrotoluene	116 a	(68 - 113)			SW846 8330
	120 a	(68 - 113)	3.0	(0-30)	SW846 8330
1,3-Dinitrobenzene	125 a	(72 - 123)			SW846 8330
	130 a	(72 - 123)	4.1	(0-29)	SW846 8330
2,4-Dinitrotoluene	120 a	(70 - 119)			SW846 8330
	124 a	(70 - 119)	2.9	(0-30)	SW846 8330
2,6-Dinitrotoluene	122 a	(71 - 119)			SW846 8330
	125 a	(71 - 119)	2.9	(0-29)	SW846 8330
HMX	118 a	(67 - 115)			SW846 8330
	120 a	(67 - 115)	1.4	(0-32)	SW846 8330
Nitrobenzene	121 a	(69 - 119)			SW846 8330
	124 a	(69 - 119)	2.5	(0-31)	SW846 8330
2-Nitrotoluene	117	(64 - 120)			SW846 8330
	113	(64 - 120)	2.7	(0-36)	SW846 8330
3-Nitrotoluene	113	(67 - 114)			SW846 8330
	115 a	(67 - 114)	2.0	(0-31)	SW846 8330
4-Nitrotoluene	114	(67 - 115)			SW846 8330
	116 a	(67 - 115)	2.3	(0-32)	SW846 8330
RDX	121	(68 - 122)			SW846 8330
	129 a	(68 - 122)	6.3	(0-32)	SW846 8330
Tetryl	102	(66 - 105)			SW846 8330
	106 a	(66 - 105)	4.2	(0-26)	SW846 8330
1,3,5-Trinitrobenzene	119	(74 - 120)			SW846 8330
	124 a	(74 - 120)	4.3	(0-29)	SW846 8330
2,4,6-Trinitrotoluene	113 a	(69 - 111)			SW846 8330
	117 a	(69 - 111)	3.4	(0-28)	SW846 8330

(Continued on next page)

MATRIX SPIKE SAMPLE EVALUATION REPORT

HPLC

Client Lot #...: G9F270189      Work Order #...: LFV9K1AF-MS      Matrix.....: WATER  
MS Lot-Sample #: G9F300242-004      LFV9K1AG-MSD

<u>SURROGATE</u>	<u>PERCENT RECOVERY</u>	<u>RECOVERY LIMITS</u>
3,4-Dinitrotoluene	104	(79 - 111)
	107	(79 - 111)

**NOTE (S) :**

Calculations are performed before rounding to avoid round-off errors in calculated results

Bold print denotes control parameters

**p** Relative percent difference (RPD) is outside stated control limits.

**a** Spiked analyte recovery is outside stated control limits.

MATRIX SPIKE SAMPLE DATA REPORT

HPLC

Client Lot #...: G9F270189      Work Order #...: LFV9K1AF-MS      Matrix.....: WATER  
 MS Lot-Sample #: G9F300242-004      LFV9K1AG-MSD  
 Date Sampled...: 06/26/09      Date Received...: 06/30/09  
 Prep Date.....: 07/01/09      Analysis Date...: 07/03/09  
 Prep Batch #...: 9182192  
 Dilution Factor: 0.98

PARAMETER	SAMPLE	SPIKE	MEASRD	UNITS	PERCNT		METHOD
	AMOUNT	AMT	AMOUNT		RECVRY	RPD	
Nitroglycerin	ND	4.91	5.19	ug/L	106		SW846 8330
	ND	4.92	5.29	ug/L	108	2.0	SW846 8330
PETN	ND	4.91	5.41	ug/L	110		SW846 8330
	ND	4.92	5.46	ug/L	111	0.90	SW846 8330
Picric Acid	ND	4.91	3.41	ug/L	70		SW846 8330
	ND	4.92	4.85	ug/L	99 p	35	SW846 8330
2-Amino-4,6-dinitrotoluene	ND	0.982	1.22	ug/L	125 a		SW846 8330
	ND	0.983	1.26	ug/L	129 a	3.4	SW846 8330
4-Amino-2,6-dinitrotoluene	ND	0.982	1.14	ug/L	116 a		SW846 8330
	ND	0.983	1.18	ug/L	120 a	3.0	SW846 8330
1,3-Dinitrobenzene	ND	0.982	1.23	ug/L	125 a		SW846 8330
	ND	0.983	1.28	ug/L	130 a	4.1	SW846 8330
2,4-Dinitrotoluene	ND	0.982	1.18	ug/L	120 a		SW846 8330
	ND	0.983	1.22	ug/L	124 a	2.9	SW846 8330
2,6-Dinitrotoluene	ND	0.982	1.19	ug/L	122 a		SW846 8330
	ND	0.983	1.23	ug/L	125 a	2.9	SW846 8330
HMX	ND	0.982	1.16	ug/L	118 a		SW846 8330
	ND	0.983	1.18	ug/L	120 a	1.4	SW846 8330
Nitrobenzene	ND	0.982	1.19	ug/L	121 a		SW846 8330
	ND	0.983	1.22	ug/L	124 a	2.5	SW846 8330
2-Nitrotoluene	ND	0.982	1.14	ug/L	117		SW846 8330
	ND	0.983	1.11	ug/L	113	2.7	SW846 8330
3-Nitrotoluene	ND	0.982	1.11	ug/L	113		SW846 8330
	ND	0.983	1.14	ug/L	115 a	2.0	SW846 8330
4-Nitrotoluene	ND	0.982	1.12	ug/L	114		SW846 8330
	ND	0.983	1.14	ug/L	116 a	2.3	SW846 8330
RDX	ND	0.982	1.19	ug/L	121		SW846 8330
	ND	0.983	1.27	ug/L	129 a	6.3	SW846 8330
Tetryl	ND	0.982	0.998	ug/L	102		SW846 8330
	ND	0.983	1.04	ug/L	106 a	4.2	SW846 8330
1,3,5-Trinitrobenzene	0.035	0.982	1.20	ug/L	119		SW846 8330
	0.035	0.983	1.26	ug/L	124 a	4.3	SW846 8330
2,4,6-Trinitrotoluene	ND	0.982	1.11	ug/L	113 a		SW846 8330
	ND	0.983	1.15	ug/L	117 a	3.4	SW846 8330

(Continued on next page)

MATRIX SPIKE SAMPLE DATA REPORT

HPLC

Client Lot #...: G9F270189      Work Order #...: LFV9K1AF-MS      Matrix.....: WATER  
MS Lot-Sample #: G9F300242-004      LFV9K1AG-MSD

<u>SURROGATE</u>	<u>PERCENT RECOVERY</u>	<u>RECOVERY LIMITS</u>
3,4-Dinitrotoluene	104	(79 - 111)
	107	(79 - 111)

**NOTE(S) :**

Calculations are performed before rounding to avoid round-off errors in calculated results.

Bold print denotes control parameters

p Relative percent difference (RPD) is outside stated control limits.

a Spiked analyte recovery is outside stated control limits.

LABORATORY CONTROL SAMPLE EVALUATION REPORT

HPLC

Client Lot #...: G9F270189      Work Order #...: LFW791AC      Matrix.....: WATER  
 LCS Lot-Sample#: G9G010000-192  
 Prep Date.....: 07/01/09      Analysis Date...: 07/02/09  
 Prep Batch #...: 9182192  
 Dilution Factor: 1

<u>PARAMETER</u>	<u>PERCENT RECOVERY</u>	<u>RECOVERY LIMITS</u>	<u>METHOD</u>
Nitroglycerin	107	(85 - 115)	SW846 8330
PETN	110	(84 - 117)	SW846 8330
Picric Acid	88	(21 - 118)	SW846 8330
2-Amino-4,6-dinitrotoluene	127 a	(77 - 123)	SW846 8330
4-Amino-2,6-dinitrotoluene	119 a	(68 - 113)	SW846 8330
1,3-Dinitrobenzene	128 a	(72 - 123)	SW846 8330
2,4-Dinitrotoluene	123 a	(70 - 119)	SW846 8330
2,6-Dinitrotoluene	125 a	(71 - 119)	SW846 8330
HMX	120 a	(67 - 115)	SW846 8330
Nitrobenzene	124 a	(69 - 119)	SW846 8330
2-Nitrotoluene	115	(64 - 120)	SW846 8330
3-Nitrotoluene	118 a	(67 - 114)	SW846 8330
4-Nitrotoluene	117 a	(67 - 115)	SW846 8330
RDX	127 a	(68 - 122)	SW846 8330
Tetryl	105	(66 - 105)	SW846 8330
1,3,5-Trinitrobenzene	127 a	(74 - 120)	SW846 8330
2,4,6-Trinitrotoluene	116 a	(69 - 111)	SW846 8330

<u>SURROGATE</u>	<u>PERCENT RECOVERY</u>	<u>RECOVERY LIMITS</u>
3,4-Dinitrotoluene	106	(79 - 111)

**NOTE(S) :**

Calculations are performed before rounding to avoid round-off errors in calculated results.

Bold print denotes control parameters

a Spiked analyte recovery is outside stated control limits

LABORATORY CONTROL SAMPLE DATA REPORT

HPLC

Client Lot #...: G9F270189      Work Order #...: LEW791AC      Matrix.....: WATER  
 LCS Lot-Sample#: G9G010000-192  
 Prep Date.....: 07/01/09      Analysis Date...: 07/02/09  
 Prep Batch #...: 9182192  
 Dilution Factor: 1

<u>PARAMETER</u>	<u>SPIKE</u> <u>AMOUNT</u>	<u>MEASURED</u> <u>AMOUNT</u>	<u>UNITS</u>	<u>PERCENT</u> <u>RECOVERY</u>	<u>METHOD</u>
Nitroglycerin	5.00	5.35	ug/L	107	SW846 8330
PETN	5.00	5.52	ug/L	110	SW846 8330
Picric Acid	5.00	4.42	ug/L	88	SW846 8330
2-Amino-4,6- dinitrotoluene	1.00	1.27 a	ug/L	127	SW846 8330
4-Amino-2,6- dinitrotoluene	1.00	1.19 a	ug/L	119	SW846 8330
1,3-Dinitrobenzene	1.00	1.28 a	ug/L	128	SW846 8330
2,4-Dinitrotoluene	1.00	1.23 a	ug/L	123	SW846 8330
2,6-Dinitrotoluene	1.00	1.25 a	ug/L	125	SW846 8330
HMX	1.00	1.20 a	ug/L	120	SW846 8330
Nitrobenzene	1.00	1.24 a	ug/L	124	SW846 8330
2-Nitrotoluene	1.00	1.15	ug/L	115	SW846 8330
3-Nitrotoluene	1.00	1.18 a	ug/L	118	SW846 8330
4-Nitrotoluene	1.00	1.17 a	ug/L	117	SW846 8330
RDX	1.00	1.27 a	ug/L	127	SW846 8330
Tetryl	1.00	1.05	ug/L	105	SW846 8330
1,3,5-Trinitrobenzene	1.00	1.27 a	ug/L	127	SW846 8330
2,4,6-Trinitrotoluene	1.00	1.16 a	ug/L	116	SW846 8330

<u>SURROGATE</u>	<u>PERCENT</u> <u>RECOVERY</u>	<u>RECOVERY</u> <u>LIMITS</u>
3,4-Dinitrotoluene	106	(79 - 111)

**NOTE(S) :**

Calculations are performed before rounding to avoid round-off errors in calculated results.

Bold print denotes control parameters

a Spiked analyte recovery is outside stated control limits.

METHOD BLANK REPORT

HPLC

Client Lot #...: G9F270189      Work Order #...: LFW791AA      Matrix.....: WATER  
 MB Lot-Sample #: G9G010000-192  
 Prep Date.....: 07/01/09  
 Analysis Date...: 07/02/09      Prep Batch #...: 9182192  
 Dilution Factor: 1

PARAMETER	RESULT	REPORTING LIMIT	UNITS	METHOD
Nitroglycerin	ND	0.65	ug/L	SW846 8330
PETN	ND	0.65	ug/L	SW846 8330
Picric Acid	ND	1.0	ug/L	SW846 8330
2-Amino-4,6-dinitrotoluene	ND	0.20	ug/L	SW846 8330
4-Amino-2,6-dinitrotoluene	ND	0.10	ug/L	SW846 8330
1,3-Dinitrobenzene	ND	0.10	ug/L	SW846 8330
2,4-Dinitrotoluene	ND	0.10	ug/L	SW846 8330
2,6-Dinitrotoluene	ND	0.10	ug/L	SW846 8330
HMX	ND	0.10	ug/L	SW846 8330
Nitrobenzene	ND	0.10	ug/L	SW846 8330
2-Nitrotoluene	ND	0.40	ug/L	SW846 8330
3-Nitrotoluene	ND	0.40	ug/L	SW846 8330
4-Nitrotoluene	ND	0.50	ug/L	SW846 8330
RDX	ND	0.10	ug/L	SW846 8330
Tetryl	ND	0.10	ug/L	SW846 8330
1,3,5-Trinitrobenzene	0.039 J	0.10	ug/L	SW846 8330
2,4,6-Trinitrotoluene	ND	0.10	ug/L	SW846 8330

SURROGATE	PERCENT RECOVERY	RECOVERY LIMITS
3,4-Dinitrotoluene	106	(79 - 111)

NOTE(S):

Calculations are performed before rounding to avoid round-off errors in calculated results.

J Estimated result. Result is less than RL.

# QC DATA ASSOCIATION SUMMARY

G9F270189

Sample Preparation and Analysis Control Numbers

<u>SAMPLE#</u>	<u>MATRIX</u>	<u>ANALYTICAL METHOD</u>	<u>LEACH BATCH #</u>	<u>PREP BATCH #</u>	<u>MS RUN#</u>
001	WATER	SW846 8330		9182192	9182194
002	WATER	SW846 8330		9182192	9182194
003	WATER	SW846 8330		9182192	9182194
004	WATER	SW846 8330		9182192	9182194
005	WATER	SW846 8330		9182192	9182194





RQC058

TestAmerica Laboratories, Inc.  
EXTRACTION BENCH WORKSHEET

Run Date: 7/02/09  
Time: 11:49:39

LEV	LEV	LEV	LEV
1	1	2	2
Y	Y	Y	Y
Y	Y	Y	Y
Y	Y	Y	Y

Blank  
Check  
MS/MSD

Weights/Volumes  
Spike & Surrogate Worksheet  
Vial contains correct volume  
Labels, greenbars, worksheets  
computer batch: correct & all match  
Anomalies to Extraction Method

Expanded Deliverable  
COC Completed  
Bench Sheet Copied  
Package Submitted to Analytical Group  
Bench Sheet Copied per COC

Extractionist: 002448 Tuan Q. Phan

Concentrationist: 002448 Tuan Q. Phan

\*\*\*\*\*  
\* QC BATCH: 9182192 \*  
\* PREP DATE: 7/01/09 9:00  
\* COMP DATE: 7/02/09 10:40  
\*\*\*\*\*

Reviewer/Date: PHANT / 7/02/09

Nitroaromatics & Nitramines: Explosives (8330)  
SOLID PHASE EXTRACTION (NOMINAL)

EXTR EXPR	ANL DUE	LOT#,MSRUN#/ WORK ORDER	TEST FLGS	EXT MTH	MATRIX	INIT/ FIN WT/VOL	PH'S ADJ1	ADJ2	EXTRACTION VOL	SOLVENTS EXCHANGE VOL	SPIKE STANDARD/ SURROGATE ID
7/01/09	7/10/09	G9F260326-001 LFOAK-1-AC		B7	A0 WATER	997.78mL 20.00mL	NA	NA	HOAC/ACN	4.5 HOAC/ACN	5.0 50UL-09GCSV0172
7/01/09	7/10/09	G9F260326-002 LFOAL-1-AC		B7	A0 WATER	991.12mL 20.00mL	NA	NA	HOAC/ACN	4.5 HOAC/ACN	5.0 50UL-09GCSV0172
7/01/09	7/10/09	G9F260326-003 LFOAM-1-AC		B7	A0 WATER	1009.82mL 20.00mL	NA	NA	HOAC/ACN	4.5 HOAC/ACN	5.0 50UL-09GCSV0172
7/01/09	7/13/09	G9F270189-001 LFRD6-1-AC		B7	A0 WATER	985.56mL 20.00mL	NA	NA	HOAC/ACN	4.5 HOAC/ACN	5.0 50UL-09GCSV0172
7/02/09	7/13/09	G9F270189-002 LFRD7-1-AC		B7	A0 WATER	999.12mL 20.00mL	NA	NA	HOAC/ACN	4.5 HOAC/ACN	5.0 50UL-09GCSV0172
7/02/09	7/13/09	G9F270189-003 LFRD8-1-AC		B7	A0 WATER	991.73mL 20.00mL	NA	NA	HOAC/ACN	4.5 HOAC/ACN	5.0 50UL-09GCSV0172
7/02/09	7/13/09	G9F270189-004 LFRD9-1-AC		B7	A0 WATER	996.78mL 20.00mL	NA	NA	HOAC/ACN	4.5 HOAC/ACN	5.0 50UL-09GCSV0172

RQC058

TestAmerica Laboratories, Inc.  
EXTRACTION BENCH WORKSHEET

Run Date: 7/02/09  
Time: 11:49:39

\*\*\*\*\*  
\* QC BATCH: 9182192 \*  
\* PREP DATE: 7/01/09 9:00  
\* COMP DATE: 7/02/09 10:40  
\*\*\*\*\*

EXTR EXPR	ANL DUE	LOT#,MSRUN#/ WORK ORDER	TEST FLGS	EXT MTH	MATRIX	INIT/ WT/VOL	PH'S ADJ1	ADJ2	EXTRACTION VOL	SOLVENTS VOL	EXCHANGE VOL	SPIKE STANDARD/ SURROGATE ID	
7/02/09	7/13/09	G9F270189-005 LFREA-1-AC		B7	A0 WATER	985.33mL 20.00mL	NA	NA	HOAC/ACN	4.5	HOAC/ACN	5.0	50UL-09GCSV0172
COMMENTS:													
7/02/09	7/14/09	G9F300242-001 LFV9A-1-AC		B7	A0 WATER	1014.4mL 20.00mL	NA	NA	HOAC/ACN	4.5	HOAC/ACN	5.0	50UL-09GCSV0172
COMMENTS:													
7/02/09	7/14/09	G9F300242-002 LFV9G-1-AC		B7	A0 WATER	975.28mL 20.00mL	NA	NA	HOAC/ACN	4.5	HOAC/ACN	5.0	50UL-09GCSV0172
COMMENTS:													
7/02/09	7/14/09	G9F300242-003 LFV9J-1-AC		B7	A0 WATER	1018.21mL 20.00mL	NA	NA	HOAC/ACN	4.5	HOAC/ACN	5.0	50UL-09GCSV0172
COMMENTS:													
7/03/09	7/14/09	G9F300242-004 LFV9K-1-AC		B7	A0 WATER	1028.11mL 20.00mL	NA	NA	HOAC/ACN	4.5	HOAC/ACN	5.0	50UL-09GCSV0172
COMMENTS:													
7/03/09	7/14/09	G9F300242-004 LFV9K-1-APS		B7	A0 WATER	1018.28mL 20.00mL	NA	NA	HOAC/ACN	4.5	HOAC/ACN	5.0	SEE BENCH SHEET 50UL-09GCSV0172
COMMENTS:													
7/03/09	7/14/09	G9F300242-004 LFV9K-1-AGD		B7	A0 WATER	1017.3mL 20.00mL	NA	NA	HOAC/ACN	4.5	HOAC/ACN	5.0	SEE BENCH SHEET 50UL-09GCSV0172
COMMENTS:													
7/03/09	7/14/09	G9F300242-005 LFV9M-1-AC		B7	A0 WATER	1016.72mL 20.00mL	NA	NA	HOAC/ACN	4.5	HOAC/ACN	5.0	50UL-09GCSV0172
COMMENTS:													
7/01/09	0/00/00	G9G010000-192 LFW79-1-AAB		B7	A0 WATER	1000mL 20.00mL	NA	NA	HOAC/ACN	4.5	HOAC/ACN	5.0	50UL-09GCSV0172
COMMENTS:													
7/01/09	0/00/00	G9G010000-192 LFW79-1-ACC		B7	A0 WATER	1000mL 20.00mL	NA	NA	HOAC/ACN	4.5	HOAC/ACN	5.0	SEE BENCH SHEET 50UL-09GCSV0172
COMMENTS:													

1% HOAC/ACN 2991-95F; 1%HOAC/H2O 2991-87B; FILTER MILLIPORE LOT R5PN29184  
SPE COLUMN WATER LOT 00333984A; SODIUM CHLORIDE MALLINCKRODT LOT H07582  
LCS, MS/MSD SPIKE 20UL-09GCSV003; 100UL-09GCSV0149; 100UL-09GCSV0039



TestAmerica West Sacramento

INITIAL CALIBRATION DATA

Start Cal Date : 10-JUN-2009 15:33  
 End Cal Date : 11-JUN-2009 13:47  
 Quant Method : ESTD  
 Origin : Disabled  
 Target Version : 4.14  
 Integrator : Falcon  
 Method file : \\Terastation\share\GCdata\LC10.I\06102009.B\8330AB.M  
 Last Edit : 11-Jun-2009 15:06 kenneyf  
 Curve Type : Average

Calibration File Names:

Level 1: \\Terastation\share\GCdata\GCdata\LC10.I\06102009.B\A-000002.d  
 Level 2: \\Terastation\share\GCdata\GCdata\LC10.I\06102009.B\A-000003.d  
 Level 3: \\Terastation\share\GCdata\GCdata\LC10.I\06102009.B\A-000015.d  
 Level 4: \\Terastation\share\GCdata\GCdata\LC10.I\06102009.B\A-000005.d  
 Level 5: \\Terastation\share\GCdata\GCdata\LC10.I\06102009.B\A-000006.d  
 Level 6: \\Terastation\share\GCdata\GCdata\LC10.I\06102009.B\A-000007.d  
 Level 7: \\Terastation\share\GCdata\GCdata\LC10.I\06102009.B\A-000008.d  
 Level 8: \\Terastation\share\GCdata\GCdata\LC10.I\06102009.B\A-000009.d

Compound	5 000	10.000	20.000	50.000	100.000	200 000	RRF	% RSD
	Level 1	Level 2	Level 3	Level 4	Level 5	Level 6		
	500.000	1000 000						
	Level 7	Level 8						
2 HMX	119 ✓ 116	120 ✓ 123	115 ✓	125 ✓	123 ✓	124	121	3.066
3 RDX	81.80000 ✓ 69.16800	75.90000 ✓ 79.81500	77.15000 ✓	80.72000 ✓	84.45000 ✓	80.13000	77.89163	5.156
4 EGDN	++++ ++++	++++ ++++	++++	++++	++++	++++	++++	++++
5 Picric ACID	++++ 74.63400	++++ 74.87800	91.72000	85.68000	84.85500	84.88200	82.77483	8.120
6 1,3,5-Trinitrobenzene	161 136	146 147	145	150	147	148	148	4 610
7 1,3-Dinitrobenzene	145 128	142 141	138	142	139	142	140	3.715
8 3,5-Dinitroaniline	97.20000 82.20800	90.80000 89.25000	89 35000	91.80000	88.83000	90.35000	89.97350	4.581
9 TETRYL	97.60000 ✓ 92.43000	95.10000 ✓ 94.40900	77.40000 ✓	96.02000 ✓	94 31000 ✓	95 06000 ✓	92.79113	6.888

*AK*  
*6/11/09*

TestAmerica West Sacramento

INITIAL CALIBRATION DATA

Start Cal Date : 10-JUN-2009 15:33  
 End Cal Date : 11-JUN-2009 13:47  
 Quant Method : ESTD  
 Origin : Disabled  
 Target Version : 4.14  
 Integrator : Falcon  
 Method file : \\Terastation\share\GCdata\LC10.I\06102009.B\8330AB.M  
 Last Edit : 11-Jun-2009 15:06 kenneyf  
 Curve Type : Average

Compound	5.000	10.000	20.000	50.000	100.000	200.000	RRF	% RSD
	Level 1	Level 2	Level 3	Level 4	Level 5	Level 6		
	500.000	1000.000						
	Level 7	Level 8						
10 Nitrobenzene	63.80000 56.02600	62.30000 61.95700	56.55000	61.44000	61.06000	62.26000	60.67412	4.659
11 Nitroglycerin	++++ ++++	++++ ++++	++++	++++	++++	++++	++++	++++
12 2,4,6-Trinitrotoluene	94.80000 84.22200	90.20000 87.17800	81.90000	88.78000	87.04000	87.55500	87.70937	4.401
13 4-AM-2,6-DNT	67.00000 58.65400	63.30000 61.14600	60.85000	63.26000	61.61000	62.67000	62.31125	3.901
14 2-AM-4,6-DNT	74.80000 69.18400	75.80000 72.13500	73.10000	75.54000	73.53000	74.39000	73.55987	2.932
15 2,6-Dinitrotoluene	57.60000 50.43000	53.30000 52.74900	51.90000	53.68000	53.03000	53.21500	53.23800	3.835
16 2,4-Dinitrotoluene	93.00000 81.43200	86.50000 85.05700	82.60000	86.20000	84.72000	85.49000	85.62488	4.031
17 2-Nitrotoluene	40.80000 33.97400	37.40000 35.45000	33.20000	35.64000	35.21000	35.77500	35.93113	6.486
18 4-Nitrotoluene	48.00000 41.67600	46.40000 43.21300	40.75000	43.56000	43.22000	43.56500	43.79800	5.390
19 3-Nitrotoluene	47.00000 41.51600	44.60000 42.82400	39.85000	43.24000	42.93000	43.08000	43.13000	4.852
20 PETN	++++ ++++	++++ ++++	++++	++++	++++	++++	++++	++++

TestAmerica West Sacramento

INITIAL CALIBRATION DATA

Start Cal Date : 10-JUN-2009 15:33  
 End Cal Date : 11-JUN-2009 13:47  
 Quant Method : ESTD  
 Origin : Disabled  
 Target Version : 4.14  
 Integrator : Falcon  
 Method file : \\Terastation\share\GCdata\LC10.I\06102009.B\8330AB.M  
 Last Edit : 11-Jun-2009 15:06 kenneyf  
 Curve Type : Average

Compound	5.000	10.000	20.000	50.000	100.000	200.000	RRF	% RSD
	Level 1	Level 2	Level 3	Level 4	Level 5	Level 6		
	500.000	1000.000						
	Level 7	Level 8						
\$ 1,3,4-Dinitrotoluene	+++++	50.10000	46.40000	46.48000	45.87000	46.85500		
	46.08000	48.38600					47.16729	3.251

TestAmerica West Sacramento

INITIAL CALIBRATION DATA

Start Cal Date : 10-JUN-2009 16:23  
 End Cal Date : 11-JUN-2009 13:47  
 Quant Method : ESTD  
 Origin : Disabled  
 Target Version : 4.14  
 Integrator : Falcon  
 Method file : \\Terastation\share\GCdata\LC10.I\06102009.B\8330AB.M\83:  
 Last Edit : 11-Jun-2009 15:14 kenneyf  
 Curve Type : Average

Calibration File Names:

Level 2: \\Terastation\share\GCdata\GCdata\LC10.I\06102009.B\A-000003.d\A-  
 Level 3: \\Terastation\share\GCdata\GCdata\LC10.I\06102009.B\A-000015.d\A-  
 Level 4: \\Terastation\share\GCdata\GCdata\LC10.I\06102009.B\A-000005.d\A-  
 Level 5: \\Terastation\share\GCdata\GCdata\LC10.I\06102009.B\A-000006.d\A-  
 Level 6: \\Terastation\share\GCdata\GCdata\LC10.I\06102009.B\A-000007.d\A-  
 Level 7: \\Terastation\share\GCdata\GCdata\LC10.I\06102009.B\A-000008.d\A-  
 Level 8: \\Terastation\share\GCdata\GCdata\LC10.I\06102009.B\A-000009.d\A-

Compound	10.000	20.000	50.000	100.000	200.000	500.000	RRF	% RSD
	Level 2	Level 3	Level 4	Level 5	Level 6	Level 7		
2 HMX	+++++	+++++	+++++	+++++	+++++	+++++	+++++	+++++
3 RDX	+++++	+++++	+++++	+++++	+++++	+++++	+++++	+++++
4 EGDN	+++++	+++++	+++++	+++++	+++++	+++++	+++++	+++++
5 Picric ACID	+++++	134	126	125	125	110	122	8.080
6 1,3,5-Trinitrobenzene	+++++	+++++	+++++	+++++	+++++	+++++	+++++	+++++
7 1,3-Dinitrobenzene	+++++	+++++	+++++	+++++	+++++	+++++	+++++	+++++
8 3,5-Dinitroaniline	+++++	+++++	+++++	+++++	+++++	+++++	+++++	+++++
9 TETRYL	+++++	+++++	+++++	+++++	+++++	+++++	+++++	+++++



TestAmerica West Sacramento

INITIAL CALIBRATION DATA

Start Cal Date : 10-JUN-2009 16:23  
 End Cal Date : 11-JUN-2009 13:47  
 Quant Method : ESTD  
 Origin : Disabled  
 Target Version : 4.14  
 Integrator : Falcon  
 Method file : \\Terastation\share\GCdata\LC10.I\06102009.B\8330AB.M\833  
 Last Edit : 11-Jun-2009 15:14 kenneyf  
 Curve Type : Average

Compound	10.000 Level 2	20.000 Level 3	50.000 Level 4	100.000 Level 5	200.000 Level 6	500.000 Level 7	RRF	% RSD
	1000.000 Level 8							
10 Nitrobenzene	+++++	+++++	+++++	+++++	+++++	+++++	+++++	+++++
11 Nitroglycerin	+++++	61.25000	59.82000	60.58000	61.86500	59.75400	60.83467	1.527
12 2,4,6-Trinitrotoluene	+++++	+++++	+++++	+++++	+++++	+++++	+++++	+++++
13 4-AM-2,6-DNT	+++++	+++++	+++++	+++++	+++++	+++++	+++++	+++++
14 2-AM-4,6-DNT	+++++	+++++	+++++	+++++	+++++	+++++	+++++	+++++
15 2,6-Dinitrotoluene	+++++	+++++	+++++	+++++	+++++	+++++	+++++	+++++
16 2,4-Dinitrotoluene	+++++	+++++	+++++	+++++	+++++	+++++	+++++	+++++
17 2-Nitrotoluene	+++++	+++++	+++++	+++++	+++++	+++++	+++++	+++++
18 4-Nitrotoluene	+++++	+++++	+++++	+++++	+++++	+++++	+++++	+++++
19 3-Nitrotoluene	+++++	+++++	+++++	+++++	+++++	+++++	+++++	+++++
20 PETN	+++++	32.65000	30.36000	31.91000	32.25000	32.38000	32.05033	2.748

TestAmerica West Sacramento

INITIAL CALIBRATION DATA

Start Cal Date : 10-JUN-2009 16:23  
 End Cal Date : 11-JUN-2009 13:47  
 Quant Method : ESTD  
 Origin : Disabled  
 Target Version : 4.14  
 Integrator : Falcon  
 Method file : \\Terastation\share\GCdata\LC10.I\06102009.B\8330AB.M\83  
 Last Edit : 11-Jun-2009 15:14 kenneyf  
 Curve Type : Average

Compound	10.000	20.000	50.000	100.000	200.000	500.000	RRF	% RSD
	Level 2	Level 3	Level 4	Level 5	Level 6	Level 7		
	1000.000							
	Level 8							
\$ 1 3,4-Dinitrotoluene	* +++++	90.40000	98.76000	88.28000	90.51000	88.03667		
	93.14200						89.85478	2.142

\* Level 2 dropped due to poor integration  
 Jky  
 6/12/09

Chromatography Summary

Injection Date: 6/10/2009 23:09 Operator: mk  
 DataFile: LC10.N06102009.BVA-000011.D Vial Num: 20  
 Instrument ID: LC10

Method 8330 Target Analyte Results

Method File: LC10.N06102009.BVA8330AB.M  
 Start Cal Date: 6/10/2009 15:33 End Cal Date: 6/11/2009 13:47

Sample: ICV\_6 08GCSV0397  
 100/200/100/100ng/mL

Dilution Factor	Extract Volume	Sample Volume	Sample Weight
1X	0 mL	0 mL	0 g

Matrix: NONE SubList: CAL.sub SpikeList:  
 Samp. Info: ICV\_6 08GCSV0397 100/200/100/100ng/mL;2  
 Misc. Info: ;6;; ;3;CAL.sub; ;0;1

Signal 1 UV 250-265										Signal 2 UV 358-205						
Compound Name	RT	Response	PPB	Spike Level	%D	Result	Flag	RT	Response	PPB	Spike Level	%D	Result	Flag	Units(%)	Flag
3,4-Dinitrotoluene				200	-100%	Fails					200	-100%	Fails		(±15)	
HMX	5.29	26153	216.4000<	200	8%	Acceptable					200	-100%	Fails		(±15)	45
RDX	7.84	16084	206.5000<	200	3%	Acceptable					200	-100%	Fails		(±15)	45
Picric ACID	9.00	41754	504.4000 ✓	500	1%	Acceptable		9.00	61500	505.5000<	500	1%	Acceptable		(±15)	
1,3,5-Trinitrobenzene	10.20	29097	197.1000<	200	-1%	Acceptable					200	-100%	Fails		(±15)	45
1,3-Dinitrobenzene	13.25	27834	199.2000<	200	0%	Acceptable					200	-100%	Fails		(±15)	45
TETRYL	14.30	17776	191.6000<	200	-4%	Acceptable					200	-100%	Fails		(±15)	45
Nitrobenzene	15.16	12143	200.1000<	200	0%	Acceptable					200	-100%	Fails		(±15)	45
2,4,6-Trinitrotoluene	16.53	17352	197.8000<	200	-1%	Acceptable					200	-100%	Fails		(±15)	45
4-AM-2,6-DNT	17.51	11969	192.1000< ✓	200	-4%	Acceptable					200	-100%	Fails		(±15)	45
2-AM-4,6-DNT	18.57	13851	188.3000<	200	-6%	Acceptable					200	-100%	Fails		(±15)	45
2,6-Dinitrotoluene	20.28	10445	196.2000<	200	-2%	Acceptable					200	-100%	Fails		(±15)	45
2,4-Dinitrotoluene	21.00	16653	194.5000<	200	-3%	Acceptable					200	-100%	Fails		(±15)	45
2-Nitrotoluene	24.68	7069	196.7000<	200	-2%	Acceptable					200	-100%	Fails		(±15)	45
4-Nitrotoluene	26.50	8526	194.7000<	200	-3%	Acceptable					200	-100%	Fails		(±15)	45
3-Nitrotoluene	28.49	8451	195.9000< ✓	200	-2%	Acceptable					200	-100%	Fails		(±15)	45
Nitroglycerin				200	-100%	Fails		15.78	12548	206.3000< ✓	200	3%	Acceptable		(±15)	45
PETN				200	-100%	Fails		30.10	5830	181.9000< ✓	200	-9%	Acceptable		(±15)	45
3,5-Dinitroaniline	13.99	18460	205.2000<	200	3%	Acceptable					200	-100%	Fails		(±15)	45
EGDN				200	-100%	Fails					200	-100%	Fails		(±15)	

Notes: M = Manually Integrated Signals Differ by More Than 40%  
 D = Operator Disabled Result Signals Differ by More Than 50%  
 O = Over Calibration Range

TestAmerica West Sacramento  
GC/LC INSTRUMENT LOG

Inst ID: LC10                      Batch ID: 07022009  
Method : Method 8330              Test : SOP SAC-LC-0009  
ICAL Date: See Calibration Report

Date	Time	Operator	Sample ID	File ID	Vol or Wt	Extract Vol	Diln	Comments
02-JUL-2009	15:27	fhk	Primer	A-000001	0 g	0 mL	1	
02-JUL-2009	16:18	fhk	Primer	A-000002	0 g	0 mL	1	
02-JUL-2009	17:09	fhk	STD_6 09GCSV0054 .4K/.2/.2/.2/	A-000003	0 g	0 mL	1	
02-JUL-2009	18:00	fhk	LFW791AA 9182192 G9G010000-192	A-000004	1000 mL	20 mL	1	
02-JUL-2009	18:50	fhk	LFW791AC 9182192 G9G010000-192	A-000005	1000 mL	20 mL	1	
02-JUL-2009	19:40	fhk	LFQAK1AC 9182192 G9F260326-1 1	A-000006	997.78 mL	20 mL	1	
02-JUL-2009	20:31	fhk	LFQAL1AC 9182192 G9F260326-2 1	A-000007	991.12 mL	20 mL	1	
02-JUL-2009	21:22	fhk	LFQAM1AC 9182192 G9F260326-3 1	A-000008	1009.82 mL	20 mL	1	
02-JUL-2009	22:13	fhk	LFRD61AC 9182192 G9F270189-1 1	A-000009	985.56 mL	20 mL	1	
02-JUL-2009	23:03	fhk	LFRD71AC 9182192 G9F270189-2 1	A-000010	999.12 mL	20 mL	1	
02-JUL-2009	23:54	fhk	LFRD81AC 9182192 G9F270189-3 1	A-000011	991.73 mL	20 mL	1	
03-JUL-2009	00:45	fhk	STD_5 09GCSV0053 .2K/.1/.1/.1/	A-000012	0 g	0 mL	1	
03-JUL-2009	01:36	fhk	LFRD91AC 9182192 G9F270189-4 1	A-000013	996.78 mL	20 mL	1	
03-JUL-2009	02:27	fhk	LFREA1AC 9182192 G9F270189-5 1	A-000014	985.33 mL	20 mL	1	
03-JUL-2009	03:17	fhk	LFV9A1AC 9182192 G9F300242-1 1	A-000015	1014.4 mL	20 mL	1	
03-JUL-2009	04:07	fhk	LFV9G1AC 9182192 G9F300242-2 1	A-000016	975.28 mL	20 mL	1	
03-JUL-2009	04:58	fhk	LFV9J1AC 9182192 G9F300242-3 1	A-000017	1018.21 mL	20 mL	1	
03-JUL-2009	05:45	fhk	LFV9K1AC 9182192 G9F300242-4 1	A-000018	1028.11 mL	20 mL	1	
03-JUL-2009	06:39	fhk	LFV9K1AF 9182192 G9F300242-4 S	A-000019	1018.28 mL	20 mL	1	
03-JUL-2009	07:30	fhk	LFV9K1AG 9182192 G9F300242-4 D	A-000020	1017.3 mL	20 mL	1	
03-JUL-2009	08:20	fhk	LFV9M1AC 9182192 G9F300242-5 1	A-000021	1016.72 mL	20 mL	1	
03-JUL-2009	09:11	fhk	STD_6 09GCSV0054 .4K/.2/.2/.2/	A-000022	0 g	0 mL	1	
03-JUL-2009	10:01	fhk	09GCSV0234 LEVEL 1	A-000023	0 g	0 mL	1	
03-JUL-2009	10:52	fhk	09GCSV0235 LEVEL 2	A-000024	0 g	0 mL	1	
03-JUL-2009	11:42	fhk	09GCSV0236 LEVEL 3	A-000025	0 g	0 mL	1	
03-JUL-2009	12:33	fhk	09GCSV0237 LEVEL 4	A-000026	0 g	0 mL	1	
03-JUL-2009	13:23	fhk	09GCSV0238 LEVEL 5	A-000027	0 g	0 mL	1	
03-JUL-2009	14:14	fhk	09GCSV0239 LEVEL 5	A-000028	0 g	0 mL	1	
03-JUL-2009	15:04	fhk	09GCSV0240 LEVEL 6	A-000029	0 g	0 mL	1	
03-JUL-2009	15:55	fhk	09GCSV0241 LEVEL 7	A-000030	0 g	0 mL	1	
03-JUL-2009	16:45	fhk	09GCSV0242 LEVEL 8	A-000031	0 g	0 mL	1	
03-JUL-2009	17:36	fhk	09GCSV0243 MRL	A-000032	0 g	0 mL	1	

Chromatography Summary

Method 8330 Target Analyte Results

Sample : **STD\_6 09GCSV0054 .4K/.2/.2/.2/.2**

Matrix: NONE SubList: CAL.sub SpikeList:  
 Samp. Info: STD\_6 09GCSV0054 .4K/ 2/.2/.2/.2  
 Misc. Info: .6. . . .3,CAL.sub, 0,1

Injection Date: 7/2/2009 17:09 Operator: fhk  
 DataFile: LC10.IA07022009.BVA-000003.D Vial Num: 2  
 Instrument ID: LC10

Method File: LC10 IA07022009.BVA8330AB.M  
 Start Cal Date: 6/10/2009 15:33 End Cal Date: 6/11/2009 13:47

Dilution Factor	Extract Volume	Sample Volume	Sample Weight
1X	0 mL	0 mL	0 g

Compound Name	RT	Signal 1 UV 250-265					Result	Flag	Signal 2 UV 358-205					Limits(%)	Flag
		Response	PPB	Spike Level	%D	Result			RT	Response	PPB	Spike Level	%D		
3,4-Dinitrotoluene	18.20	9636	204.3000<	200	7%	Acceptable		18.20	18445	213.4000	200	7%	Acceptable	(±15)	
HMX	5.29	25432	210.5000<	200	5%	Acceptable					200	-100%	Fails	(±15)	45
RDX	7.85	16295	209.2000<	200	5%	Acceptable					200	-100%	Fails	(±15)	45
Picric ACID	8.99	41811	505.1000	500	1%	Acceptable		8.99	61520	505.6000<	500	1%	Acceptable	(±15)	
1,3,5-Trinitrobenzene	10.20	30367	205.7000<	200	3%	Acceptable					200	-100%	Fails	(±15)	45
1,3-Dinitrobenzene	13.26	29099	208.3000<	200	4%	Acceptable					200	-100%	Fails	(±15)	45
TETRYL	14.32	19114	206.0000<	200	3%	Acceptable					200	-100%	Fails	(±15)	45
Nitrobenzene	15.19	12872	212.1000<	200	6%	Acceptable					200	-100%	Fails	(±15)	45
2,4,6-Trinitrotoluene	16.55	18075	206.1000<	200	3%	Acceptable					200	-100%	Fails	(±15)	45
4-AM-2,6-DNT	17.55	12873	206.6000<	200	3%	Acceptable					200	-100%	Fails	(±15)	45
2-AM-4,6-DNT	18.61	15046	204.5000<	200	2%	Acceptable					200	-100%	Fails	(±15)	45
2,6-Dinitrotoluene	20.34	10970	206.0000<	200	3%	Acceptable					200	-100%	Fails	(±15)	45
2,4-Dinitrotoluene	21.06	17567	205.2000<	200	3%	Acceptable					200	-100%	Fails	(±15)	45
2-Nitrotoluene	24.79	7368	205.0000<	200	3%	Acceptable					200	-100%	Fails	(±15)	45
4-Nitrotoluene	26.61	8982	205.1000<	200	3%	Acceptable					200	-100%	Fails	(±15)	45
3-Nitrotoluene	28.64	8962	207.8000<	200	4%	Acceptable					200	-100%	Fails	(±15)	45
Nitroglycerin				200	-100%	Fails		15.81	12531	206.0000<	200	3%	Acceptable	(±15)	45
PETN				200	-100%	Fails		30.22	6862	214.1000<	200	7%	Acceptable	(±15)	45
3,5-Dinitroaniline	14.01	18502	205.6000<	200	3%	Acceptable					200	-100%	Fails	(±15)	45
EGDN				200	-100%	Fails					200	-100%	Fails	(±15)	

Notes: M = Manually Integrated Signals Differ by More Than 40%  
 D = Operator Disabled Result Signals Differ by More Than 50%  
 O = Over Calibration Range

Chromatography Summary

Injection Date: 7/3/2009 0:45 Operator: fhk  
 DataFile: LC10 R07022009.BVA-000012.D Vial Num: 3  
 Instrument ID: LC10

Method 8330 Target Analyte Results

Sample: STD\_5 09GCSV0053.2K/1/1/1/1

Method File: LC10 R07022009.BV8330AB.M  
 Start Cal Date: 6/10/2009 15:33 End Cal Date: 6/11/2009 13:47

Matrix: NONE SubList: CAL.sub SpikeList:  
 Samp. Info: STD\_5 09GCSV0053.2K/1/1/1/1;2  
 Misc. Info: ;5; ; ;3;CAL.sub ;0,1

Dilution Factor	Extract Volume	Sample Volume	Sample Weight
1X	0 mL	0 mL	0 g

Signal 1 UV 250-265							Signal 2 UV 358-205									
Compound Name	RT	Response	PPB	Spike Level	%D	Result	Flag	RT	Response	PPB	Spike Level	%D	Result	Flag	Limits(%)	Flag
3,4-Dinitrotoluene	18.21	4770	101.1000<	100	1%	Acceptable		18.21	9105	105.3000	100	5%	Acceptable		(±15)	
HMX	5.29	12664	104.8000<	100	5%	Acceptable					100	-100%	Fails		(±15)	45
RDX	7.86	8028	103.1000<	100	3%	Acceptable					100	-100%	Fails		(±15)	45
Picric ACID	9.05	16980	205.1000	200	3%	Acceptable		9.05	25011	205.5000<	200	3%	Acceptable		(±15)	
1,3,5-Trinitrobenzene	10.21	15116	102.4000<	100	2%	Acceptable					100	-100%	Fails		(±15)	45
1,3-Dinitrobenzene	13.27	14453	103.5000<	100	4%	Acceptable					100	-100%	Fails		(±15)	45
TETRYL	14.34	9510	102.5000<	100	3%	Acceptable					100	-100%	Fails		(±15)	45
Nitrobenzene	15.20	6322	104.2000<	100	4%	Acceptable					100	-100%	Fails		(±15)	45
2,4,6-Trinitrotoluene	16.57	9018	102.8000<	100	3%	Acceptable					100	-100%	Fails		(±15)	45
4-AM-2,6-DNT	17.56	6403	102.8000<	100	3%	Acceptable					100	-100%	Fails		(±15)	45
2-AM-4,6-DNT	18.63	7508	102.1000<	100	2%	Acceptable					100	-100%	Fails		(±15)	45
2,6-Dinitrotoluene	20.35	5455	102.5000<	100	3%	Acceptable					100	-100%	Fails		(±15)	45
2,4-Dinitrotoluene	21.08	8726	101.9000<	100	2%	Acceptable					100	-100%	Fails		(±15)	45
2-Nitrotoluene	24.79	3635	101.2000<	100	1%	Acceptable					100	-100%	Fails		(±15)	45
4-Nitrotoluene	26.61	4448	101.6000<	100	2%	Acceptable					100	-100%	Fails		(±15)	45
3-Nitrotoluene	28.63	4394	101.9000<	100	2%	Acceptable					100	-100%	Fails		(±15)	45
Nitroglycerin				100	-100%	Fails		15.82	6254	102.8000<	100	3%	Acceptable		(±15)	45
PETN				100	-100%	Fails		30.25	3304	103.1000<	100	3%	Acceptable		(±15)	45
3,5-Dinitroaniline	14.02	9179	102.0000<	100	2%	Acceptable					100	-100%	Fails		(±15)	45
EGDN				100	-100%	Fails					100	-100%	Fails		(±15)	45



Notes M = Manually Integrated Signals Differ by More Than 40%  
 D = Operator Disabled Result Signals Differ by More Than 50%  
 O = Over Calibration Range

Chromatography Summary

Method 8330 Target Analyte Results

Sample: **STD\_6 09GCSV0054 .4K/.2/.2/.2**

Matrix: NONE SubList: CAL sub SpikeList:  
 Samp. Info: STD\_6 09GCSV0054 4K/2/2/2/2  
 Misc. Info: ;6, ; ; 3;CAL sub, ;0,1

Injection Date: 7/3/2009 9:11 Operator: fhk  
 DataFile: I:\C10\107022009\BVA-000022.D Vial Num: 4  
 Instrument ID: I.C10

Method File: I:\C10\107022009\B\8330AB.M  
 Start Cal Date: 6/10/2009 15:33 End Cal Date: 6/11/2009 13:47

Dilution Factor	Extract Volume	Sample Volume	Sample Weight
1X	0 mL	0 mL	0 g

Signal 1 UV 250-265								Signal 2 UV 358-205								
Compound Name	RT	Response	PPB	Spike Level	%D	Result	Flag	RT	Response	PPB	Spike Level	%D	Result	Flag	Limits(%)	Flag
3,4-Dinitrotoluene	18.24	9589	203.3000<	200	2%	Acceptable		18.23	18290	211.6000	200	6%	Acceptable		(±15)	
HMX	5.30	25440	210.6000<	200	5%	Acceptable					200	-100%	Fails		(±15)	45
RDX	7.87	16268	208.8000<	200	4%	Acceptable					200	-100%	Fails		(±15)	45
Picric ACID	9.04	42655	515.3000	500	3%	Acceptable		9.04	62723	515.5000<	500	3%	Acceptable		(±15)	
1,3,5-Trinitrobenzene	10.22	30282	205.1000<	200	3%	Acceptable					200	-100%	Fails		(±15)	45
1,3-Dinitrobenzene	13.29	28991	207.5000<	200	4%	Acceptable					200	-100%	Fails		(±15)	45
TETRYL	14.35	19191	206.8000<	200	3%	Acceptable					200	-100%	Fails		(±15)	45
Nitrobenzene	15.21	12738	209.9000<	200	5%	Acceptable					200	-100%	Fails		(±15)	45
2,4,6-Trinitrotoluene	16.58	18022	205.5000<	200	3%	Acceptable					200	-100%	Fails		(±15)	45
4-AM-2,6-DNT	17.58	12828	205.9000<	200	3%	Acceptable					200	-100%	Fails		(±15)	45
2-AM-4,6-DNT	18.65	15059	204.7000<	200	2%	Acceptable					200	-100%	Fails		(±15)	45
2,6-Dinitrotoluene	20.38	10882	204.4000<	200	2%	Acceptable					200	-100%	Fails		(±15)	45
2,4-Dinitrotoluene	21.10	17483	204.2000<	200	2%	Acceptable					200	-100%	Fails		(±15)	45
2-Nitrotoluene	24.82	7294	203.0000<	200	2%	Acceptable					200	-100%	Fails		(±15)	45
4-Nitrotoluene	26.64	8876	202.6000<	200	1%	Acceptable					200	-100%	Fails		(±15)	45
3-Nitrotoluene	28.67	8836	204.9000<	200	2%	Acceptable					200	-100%	Fails		(±15)	45
Nitroglycerin				200	-100%	Fails		15.84	12481	205.2000<	200	3%	Acceptable		(±15)	45
PETN				200	-100%	Fails		30.27	6681	208.4000<	200	4%	Acceptable		(±15)	45
3,5-Dinitroaniline	14.03	18458	205.1000<	200	3%	Acceptable					200	-100%	Fails		(±15)	45
EGDN				200	-100%	Fails					200	-100%	Fails		(±15)	



Notes: M = Manually Integrated Signals Differ by More Than 40%  
 D = Operator Disabled Result Signals Differ by More Than 50%  
 O = Over Calibration Range

23LCMW03DW

TestAmerica West Sacramento

A-000009.D

Chromatography Summary

Injection Date: 7/2/2009 22:13 Operator: rhk  
Data File: LC10 107022009 BVA-000009.D Vial Num: 16  
Instrument ID: LC10

Method 8330 Target Analyte Results

Sample: LFRD61AC 9182192 G9F270189-1 1X

Method File: LC10.107022009.BV8330AB.M  
Start Cal Date: 6/10/2009 15:33 End Cal Date: 6/11/2009 13:47

Matrix: WATER SubList: WATER sub SpikeList:  
Samp. Info: LFRD61AC 9182192 G9F270189-1 1X,0  
Misc. Info: .,985.56,,20,1,WATER sub,,0,1,LFRD61AC

Dilution Factor	Extract Volume	Sample Volume	Sample Weight
1X	20 mL	985.56 mL	0 g

Signal 1 UV 250-265						Signal 2 UV 358-205							
Compound Name	RT	Diff	Response	Conc (ug/L)	Flag	RT	Diff	Response	Conc (ug/L)	Flag	MDL	RL	Flag
3,4-Dinitrotoluene	18.26	0.020	6239	2.6800<		18.25	0.014	11996	2.8160		0.0000	0.00	
HMX											0.0274	0.10	
RDX											0.0660	0.10	
Picric ACID											0.1268	1.01	
1,3,5-Trinitrobenzene	10.14	-0.090	256	0.0352<							0.0315	0.10	45
1,3-Dinitrobenzene											0.0507	0.10	
TETRYL											0.0507	0.10	
Nitrobenzene											0.0507	0.10	
2,4,6-Trinitrotoluene											0.0244	0.10	
4-AM-2,6-DNT											0.0223	0.10	
2-AM-4,6-DNT											0.1015	0.20	
2,6-Dinitrotoluene											0.0507	0.10	
2,4-Dinitrotoluene											0.0507	0.10	
2-Nitrotoluene											0.0731	0.51	
4-Nitrotoluene											0.0731	0.51	
3-Nitrotoluene											0.0629	0.51	
Nitroglycerin											0.3348	0.66	
PETN											0.3044	0.66	
3,5-Dinitroaniline											0.0254	1.01	

ND

Surrogates:	Spiked	Recovered	% Rec	Spiked	Recovered	% Rec	Limits
3,4-Dinitrotoluene	2.5366	2.6800	106	2.5366	2.8160	111	(79-111)

Notes: M = Manually Integrated Signals Differ by More Than 40%  
D = Operator Disabled Result Signals Differ by More Than 50%  
O = Over Calibration Range



23LCMW03SW

A-000010.D

Chromatography Summary

Injection Date: 7/2/2009 23:03 Operator: fhk  
 Data File: LC10 I07022009 BVA-000010.D Vial Num: 17  
 Instrument ID: LC10

Method 8330 Target Analyte Results

Sample: LFRD71AC 9182192 G9F270189-2 1X

Method File: LC10 I07022009 BV8330AB.M  
 Start Cal Date: 6/10/2009 15:33 End Cal Date: 6/11/2009 13:47

Matrix: WATER SubList: WATER.sub SpikeList:  
 Samp. Info: LFRD71AC 9182192 G9F270189-2 1X,0  
 Misc. Info: .,999 12;.20,1,WATER sub;.0,1,LFRD71AC

Dilution Factor	Extract Volume	Sample Volume	Sample Weight
1X	20 mL	999.12 mL	0 g

Signal 1 UV 250-265						Signal 2 UV 358-205							
Compound Name	RT	Diff	Response	Conc (ug/L)	Flag	RT	Diff	Response	Conc (ug/L)	Flag	MDL	RL	Flag
3,4-Dinitrotoluene	18.25	0.017	6242	2.6490<		18.25	0.010	11999	2.7780		0.0000	0.00	
HMX											0.0270	0.10	
RDX											0.0651	0.10	
Picric ACID											0.1251	1.00	
1,3,5-Trinitrobenzene	10.12	-0.103	275	0.0373<							0.0310	0.10	45
1,3-Dinitrobenzene											0.0500	0.10	
TETRYL											0.0500	0.10	
Nitrobenzene											0.0500	0.10	
2,4,6-Trinitrotoluene											0.0240	0.10	
4-AM-2,6-DNT											0.0220	0.10	
2-AM-4,6-DNT											0.1001	0.20	
2,6-Dinitrotoluene											0.0500	0.10	
2,4-Dinitrotoluene											0.0500	0.10	
2-Nitrotoluene											0.0721	0.50	
4-Nitrotoluene											0.0721	0.50	
3-Nitrotoluene											0.0621	0.50	
Nitroglycerin											0.3303	0.65	
PETN											0.3003	0.65	
3,5-Dinitroaniline											0.0250	1.00	

ND

Surrogates	Spiked	Recovered	% Rec	Spiked	Recovered	% Rec	Limits
3,4-Dinitrotoluene	2.5022	2.6490	106	2.5022	2.7780	111	(79-111)

Notes: M = Manually Integrated Signals Differ by More Than 40%  
 D = Operator Disabled Result Signals Differ by More Than 50%  
 O = Over Calibration Range

23LC MWO20W

A-000011.D

Chromatography Summary

Injection Date: 7/2/2009 23:54 Operator: fbk
DataFile: LC10 I07022009 BAA-000011.D Vial Num: 18
Instrument ID: LC10

Method 8330 Target Analyte Results

Sample: LFRD81AC 9182192 G9F270189-3 1X

Method File: LC10 I07022009 BA8330AB.M
Start Cal Date: 6/10/2009 15:33 End Cal Date: 6/11/2009 13:47

Matrix: WATER SubList: WATER.sub SpikeList:
Samp. Info: LFRD81AC 9182192 G9F270189-3 1X;0;
Misc. Info: ,.991.73;;20,1,WATER sub,;0 1,LFRD81AC

Table with 4 columns: Dilution Factor, Extract Volume, Sample Volume, Sample Weight. Values: 1X, 20 mL, 991.73 mL, 0 g

Main chromatography table with columns for Compound Name, RT, Diff, Response, Conc (ug/L), Flag, MDL, RL, and Flag. Includes handwritten note 'not confirmed' and 'U'.

Summary table with columns: Surrogates, Spiked, Recovered, % Rec, Limits. Row for 3,4-Dinitrotoluene showing 2.5208 spiked and 2.6320 recovered.

Notes: M = Manually Integrated Signals Differ by More Than 40%
D = Operator Disabled Result Signals Differ by More Than 50%
O = Over Calibration Range

236CMW02SW

TestAmerica West Sacramento

A-000013.D

Chromatography Summary

Injection Date: 7/3/2009 1:36 Operator: fnk  
Data File: LC10 107022009 BIA-000013.D Vial Num: 19  
Instrument ID: LC10

Method 8330 Target Analyte Results

Method File: LC10 107022009.B\8330AB.M  
Start Cal Date: 6/10/2009 15:33 End Cal Date: 6/11/2009 13:47

Sample: LFRD91AC 9182192 G9F270189-4 1X

Dilution Factor	Extract Volume	Sample Volume	Sample Weight
1X	20 mL	996.78 mL	0 g

Matrix: WATER SubList: WATER sub SpikeList:  
Samp. Info: LFRD91AC 9182192 G9F270189-4 1X,0,  
Misc. Info: ;,996.78;20,1;WATER sub;;0,1;LFRD91AC

Signal 1 UV 250-265						Signal 2 UV 358-205							
Compound Name	RT	Diff	Response	Conc (ug/L)	Flag	RT	Diff	Response	Conc (ug/L)	Flag	MDL	RL	Flag
3,4-Dinitrotoluene	18.23	0.017	6192	2.6340<		18.23	-0.010	11918	2.7660		0.0000	0.00	
HMX											0.0271	0.10	
RDX											0.0652	0.10	
Picric ACID											0.1254	1.00	
1,3,5-Trinitrobenzene	10.13	-0.086	271	0.0368<							0.0311	0.10	45
1,3-Dinitrobenzene											0.0502	0.10	
TETRYL											0.0502	0.10	
Nitrobenzene											0.0502	0.10	
2,4,6-Trinitrotoluene											0.0241	0.10	
4-AM-2,6-DNT											0.0221	0.10	
2-AM-4,6-DNT											0.1003	0.20	
2,6-Dinitrotoluene											0.0502	0.10	
2,4-Dinitrotoluene											0.0502	0.10	
2-Nitrotoluene											0.0722	0.50	
4-Nitrotoluene											0.0722	0.50	
3-Nitrotoluene											0.0622	0.50	
Nitroglycenn											0.3311	0.65	
PETN											0.3010	0.65	
3,5-Dinitroaniline											0.0251	1.00	

ND

Surrogates:	Spiked	Recovered	% Rec	Spiked	Recovered	% Rec	Limits
3,4-Dinitrotoluene	2.5081	2.6340	105	2.5081	2.7660	110	(79-111)

Notes: M = Manually Integrated Signals Differ by More Than 40%  
D = Operator Disabled Result Signals Differ by More Than 50%  
O = Over Calibration Range

23 LCMWD 10W

TestAmerica West Sacramento

A-000014.D

Chromatography Summary

Injection Date: 7/3/2009 2:27 Operator: fhk  
Data File: LC10 107022009 BVA-000014.D Vial Num: 20  
Instrument ID: LC10

Method 8330 Target Analyte Results

Sample: LFREA1AC 9182192 G9F270189-5 1X

Method File: LC10 107022009 BVA8330AB.M  
Start Cal Date: 6/10/2009 15:33 End Cal Date: 6/11/2009 13:47

Matrix: WATER SubList: WATER.sub SpikeList:  
Samp. Info: LFREA1AC 9182192 G9F270189-5 1X,0,  
Misc. Info: ,,985.33,,20.1,WATER.sub,,0;1:LFREA1AC

Dilution Factor	Extract Volume	Sample Volume	Sample Weight
1X	20 mL	985.33 mL	0 g

Signal 1 UV 250-265						Signal 2 UV 358-205							
Compound Name	RT	Diff	Response	Conc (ug/L)	Flag	RT	Diff	Response	Conc (ug/L)	Flag	MDL	RL	Flag
3,4-Dinitrotoluene	18.25	0.034	6173	2.6560<		18.24	0.003	11870	2.7870		0.0000	0.00	
HMX											0.0274	0.10	
RDX											0.0660	0.10	
Picric ACID											0.1269	1.01	
1,3,5-Trinitrobenzene	10.13	-0.086	274	0.0377<							0.0315	0.10	45
1,3-Dinitrobenzene											0.0507	0.10	
1EBRYL											0.0507	0.10	
Nitrobenzene											0.0507	0.10	
2,4,6-Trinitrotoluene											0.0244	0.10	
4-AM-2,6-DNT											0.0223	0.10	
2-AM-4,6-DNT											0.1015	0.20	
2,6-Dinitrotoluene											0.0507	0.10	
2,4-Dinitrotoluene											0.0507	0.10	
2-Nitrotoluene											0.0731	0.51	
4-Nitrotoluene											0.0731	0.51	
3-Nitrotoluene											0.0629	0.51	
Nitroglycerin											0.3349	0.66	
PETN						30.09	-0.166	536	0.3394<		0.3045	0.66	45
3,5-Dinitroaniline											0.0254	1.01	

ND

MD  
7/7/09

Surrogates	Spiked	Recovered	% Rec	Spiked	Recovered	% Rec	Limits
3,4-Dinitrotoluene	2.5372	2.6560	105	2.5372	2.7870	110	(79-111)

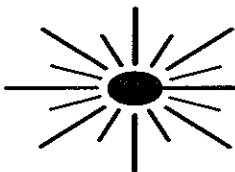
Notes: M = Manually Integrated Signals Differ by More Than 40%  
D = Operator Disabled Result Signals Differ by More Than 50%  
O = Over Calibration Range

**Data Validation Report**

**Michael Baker Jr., Inc**

**Camp Bonneville**

**SDG#: PSF0891**

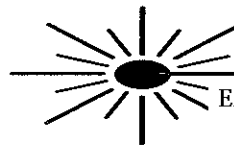


**ECT.CON INC.**

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Environmental and Computer  
Technology Consultants

3531 Fox Chase Drive  
Imperial, PA 15126  
(724) 695-8042  
FAX (724) 695-2698  
e-mail: [ectconinc@comcast.net](mailto:ectconinc@comcast.net)



## Data Validation Report

SDG#	PSF0891
Validation Report Date	August 17, 2009
Validation Guidance	USEPA CLP National Functional Guidelines for Data Review
Client Name	Michael Baker Jr., Inc.
Project Name	Camp Bonneville
Laboratory	TestAmerica
Analytical Parameters	Explosives
Analytical Method	SW-846 8330

### Samples/Matrix:

Date Sampled	Location	Portland ID	Sacramento ID	Explosives	Matrix
06/25/09	23LCMW01SW	PSF0891-01	LFV9A	X	Aqueous
06/25/09	23LF4MW17SW	PSF0891-02	LFV9G	X	Aqueous
06/25/09	23LF4MW18SW	PSF0891-03	LFV9J	X	Aqueous
06/26/09	23LF4MW7B	PSF0891-04	LFV9K	X	Aqueous
06/26/09	23LF4MW5A	PSF0891-05	LFV9M	X	Aqueous

Analytical data in this report were screened to determine analytical limitations of the data based on specific quality control criteria. This screening assumes analytical results are correct as reported and merely provides an interpretation of the reported quality control results. Laboratory calculations have been verified as part of this validation. Specific findings on analytical limitations are presented in this report. Annotated Form 1s or spreadsheets for samples reviewed are included after the Data Assessment Findings. Form 1s for the MS/MSD samples and spreadsheets are not annotated.

### SUMMARY

The sample set for the Camp Bonneville site consists of five aqueous field samples. These samples were analyzed for explosives as provided in the Table 1. The findings presented in this review of the analytical data assume that the information presented by the analytical laboratory is correct.

The explosives findings are based upon the assessment of the following:

- \* ● Data Completeness
- \* ● Holding Times
- \* ● Calibration (Initial and Continuing)
- Blanks
- \* ● System Monitoring Compounds (Surrogate Spikes)
- \* ● Matrix Spike/Matrix Spike Duplicates
- Laboratory Control Standard Results
- \* ● Target Compound Identification
- \* ● Compound Quantification and Reported Contract Quantitation Limits

- \* • System Performance
- \* Criteria were met for this evaluation item.

This evaluation was conducted in accordance with USEPA CLP National Functional Guidelines for Organic Data Review and the analytical method. Findings from this evaluation should be considered when using the analytical data. This report presents a summary of the data qualifications based on the review of the aforementioned evaluation criteria. This is followed by annotated Form 1s/ spreadsheets. Finally, the worksheets used to perform the evaluation are provided.

## FINDINGS

### 1. Blanks

The laboratory method blank exhibited contamination for the following parameter:

Blank	Compound	Maximum Concentration (ppb)	Action Limit (ppm)	Action
LFW791	1,3,5-Trinitrobenzene	0.039	0.195	U sample results < RL

RL - reporting limit

### 2. Laboratory Control Sample Results

Recovery of HMX (120%) and RDX (127%) exceeded the laboratory's upper quality control limits. In the following sample, positive results for HMX and RDX were qualified as estimated "J".

23LF4MW5A

## NOTES

### Matrix Spike/Matrix Spike Duplicate Results


A MS/MSD was performed on sample 23LF4MW7B. Recovery of 13 out of 17 compounds exceeded the laboratory's upper quality control limits. The affected compounds were not detected in the unspiked sample. Data were not qualified on this basis.

### Laboratory Control Sample Results

Recovery of 12 out of 17 compounds exceeded the laboratory's upper quality control limits. The affected compounds were not detected in the associated samples. Data were not qualified on this basis.

### Field Duplicate Results

No field duplicates were included with this SDG. Data were not qualified on this basis.

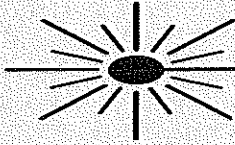
  
Data Reviewer

8/17/09  
Date

### **Glossary of Data Qualifiers**

U	Not Detected.	The associated number indicates approximate sample concentration necessary to be detected.
UJ	Not Detected.	Quantitation limit may be inaccurate or imprecise.
J	Analyte Present.	Reported value may not be accurate or precise.
N	Consider Present.	Tentative identification. Special methods may be needed to confirm its presence or absence in future sampling efforts.
R	Unusable Result.	Analyte may or may not be present in the sample.
UR	Unusable Result.	Analyte may or may not be present in the sample.





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*Annotated Form 1's*  
*(Spreadsheet)*

TestAmerica Portland

Client Sample ID: PSF0891-01

HPLC

23K MW015W

Lot-Sample #...: G9F300242-001    Work Order #...: LFV9A1AC    Matrix.....: WATER  
 Date Sampled...: 06/25/09    Date Received...: 06/30/09  
 Prep Date.....: 07/01/09    Analysis Date...: 07/03/09  
 Prep Batch #...: 9182192  
 Dilution Factor: 0.98    Method.....: SW846 8330

PARAMETER	RESULT	REPORTING LIMIT	UNITS	MDL
Nitroglycerin	ND	0.64	ug/L	0.15
PETN	ND	0.64	ug/L	0.23
Picric Acid	ND	0.98	ug/L	0.12
2-Amino-4,6-dinitrotoluene	ND	0.20	ug/L	0.098
4-Amino-2,6-dinitrotoluene	ND	0.098	ug/L	0.022
1,3-Dinitrobenzene	ND	0.098	ug/L	0.049
2,4-Dinitrotoluene	ND	0.098	ug/L	0.049
2,6-Dinitrotoluene	ND	0.098	ug/L	0.049
HMX	ND	0.098	ug/L	0.026
Nitrobenzene	ND	0.098	ug/L	0.049
2-Nitrotoluene	ND	0.39	ug/L	0.071
3-Nitrotoluene	ND	0.39	ug/L	0.061
4-Nitrotoluene	ND	0.49	ug/L	0.071
RDX	ND	0.098	ug/L	0.064
Tetryl	ND	0.098	ug/L	0.049
1,3,5-Trinitrobenzene	<del>0.032</del> -J,B	0.098	ug/L	0.030
2,4,6-Trinitrotoluene	ND	0.098	ug/L	0.024
	PERCENT	RECOVERY		
<u>SURROGATE</u>	<u>RECOVERY</u>	<u>LIMITS</u>		
3,4-Dinitrotoluene	104	(79 - 111)		

NOTE(S):

J Estimated result. Result is less than RL.

B Method blank contamination. The associated method blank contains the target analyte at a reportable level.

OK  
8/6/09

TestAmerica Portland

Client Sample ID: PSF0891-02

HPLC

23LF4.MW178W

Lot-Sample #...: G9F300242-002 Work Order #...: LFV9G1AC Matrix.....: WATER  
 Date Sampled...: 06/25/09 Date Received...: 06/30/09  
 Prep Date.....: 07/01/09 Analysis Date...: 07/03/09  
 Prep Batch #...: 9182192  
 Dilution Factor: 1.02 Method.....: SW846 8330

PARAMETER	RESULT	REPORTING		
		LIMIT	UNITS	MDL
Nitroglycerin	ND	0.66	ug/L	0.15
PETN	ND	0.66	ug/L	0.23
Picric Acid	ND	1.0	ug/L	0.13
2-Amino-4,6-dinitrotoluene	ND	0.20	ug/L	0.10
4-Amino-2,6-dinitrotoluene	ND	0.10	ug/L	0.022
1,3-Dinitrobenzene	ND	0.10	ug/L	0.051
2,4-Dinitrotoluene	ND	0.10	ug/L	0.051
2,6-Dinitrotoluene	ND	0.10	ug/L	0.051
HMX	ND	0.10	ug/L	0.028
Nitrobenzene	ND	0.10	ug/L	0.051
2-Nitrotoluene	ND	0.41	ug/L	0.073
3-Nitrotoluene	ND	0.41	ug/L	0.063
4-Nitrotoluene	ND	0.51	ug/L	0.073
RDX	ND	0.10	ug/L	0.066
Tetryl	ND	0.10	ug/L	0.051
1,3,5-Trinitrobenzene	<del>0.035 J,B</del>	0.10	ug/L	0.032
2,4,6-Trinitrotoluene	ND	0.10	ug/L	0.024
	PERCENT	RECOVERY		
SURROGATE	RECOVERY	LIMITS		
3,4-Dinitrotoluene	105	(79 - 111)		

NOTE(S):

J Estimated result. Result is less than RL.

B Method blank contamination. The associated method blank contains the target analyte at a reportable level.

WJ  
8/16/09

TestAmerica Portland

Client Sample ID: PSF0891-03

HPLC

23LF4 MW 185W

Lot-Sample #...: G9F300242-003    Work Order #...: Lfv9J1AC    Matrix.....: WATER  
 Date Sampled...: 06/25/09    Date Received...: 06/30/09  
 Prep Date.....: 07/01/09    Analysis Date...: 07/03/09  
 Prep Batch #...: 9182192  
 Dilution Factor: 0.98    Method.....: SW846 8330

PARAMETER	RESULT	REPORTING		
		LIMIT	UNITS	MDL
Nitroglycerin	ND	0.64	ug/L	0.15
PETN	ND	0.64	ug/L	0.23
Picric Acid	ND	0.98	ug/L	0.12
2-Amino-4,6-dinitrotoluene	ND	0.20	ug/L	0.098
4-Amino-2,6-dinitrotoluene	ND	0.098	ug/L	0.022
1,3-Dinitrobenzene	ND	0.098	ug/L	0.049
2,4-Dinitrotoluene	ND	0.098	ug/L	0.049
2,6-Dinitrotoluene	ND	0.098	ug/L	0.049
HMX	ND	0.098	ug/L	0.026
Nitrobenzene	ND	0.098	ug/L	0.049
2-Nitrotoluene	ND	0.39	ug/L	0.071
3-Nitrotoluene	ND	0.39	ug/L	0.061
4-Nitrotoluene	ND	0.49	ug/L	0.071
RDX	ND	0.098	ug/L	0.064
Tetryl	ND	0.098	ug/L	0.049
1,3,5-Trinitrobenzene	<del>0.037-J,B</del>	0.098 U	ug/L	0.030
2,4,6-Trinitrotoluene	ND	0.098	ug/L	0.024
	PERCENT	RECOVERY		
<u>SURROGATE</u>	<u>RECOVERY</u>	<u>LIMITS</u>		
3,4-Dinitrotoluene	103	(79 - 111)		

NOTE(S):

- J Estimated result Result is less than RL.
- B Method blank contamination. The associated method blank contains the target analyte at a reportable level

UK  
5/6/09

TestAmerica Portland

Client Sample ID: PSF0891-04

HPLC

23154 MW7B

Lot-Sample #...: G9F300242-004    Work Order #...: LFV9K1AC    Matrix.....: WATER  
 Date Sampled...: 06/26/09    Date Received...: 06/30/09  
 Prep Date.....: 07/01/09    Analysis Date...: 07/03/09  
 Prep Batch #...: 9182192  
 Dilution Factor: 0.97    Method.....: SW846 8330

PARAMETER	RESULT	REPORTING		
		LIMIT	UNITS	MDL
Nitroglycerin	ND	0.63	ug/L	0.15
PETN	ND	0.63	ug/L	0.22
Picric Acid	ND	0.97	ug/L	0.12
2-Amino-4,6-dinitrotoluene	ND	0.19	ug/L	0.097
4-Amino-2,6-dinitrotoluene	ND	0.097	ug/L	0.021
1,3-Dinitrobenzene	ND	0.097	ug/L	0.048
2,4-Dinitrotoluene	ND	0.097	ug/L	0.048
2,6-Dinitrotoluene	ND	0.097	ug/L	0.048
HMX	ND	0.097	ug/L	0.026
Nitrobenzene	ND	0.097	ug/L	0.048
2-Nitrotoluene	ND	0.39	ug/L	0.070
3-Nitrotoluene	ND	0.39	ug/L	0.060
4-Nitrotoluene	ND	0.48	ug/L	0.070
RDX	ND	0.097	ug/L	0.063
Tetryl	ND	0.097	ug/L	0.048
1,3,5-Trinitrobenzene	<del>0.035</del> J,B	0.097 U	ug/L	0.030
2,4,6-Trinitrotoluene	ND	0.097	ug/L	0.023
	PERCENT	RECOVERY		
SURROGATE	RECOVERY	LIMITS		
3,4-Dinitrotoluene	98	(79 - 111)		

NOTE(S):

J Estimated result. Result is less than RL.

B Method blank contamination. The associated method blank contains the target analyte at a reportable level.

WJ  
8/16/09

TestAmerica Portland

Client Sample ID: PSF0891-05

HPLC *23LF4MWSA*

Lot-Sample #...: G9F300242-005 Work Order #...: LFV9M1AC Matrix.....: WATER  
 Date Sampled...: 06/26/09 Date Received...: 06/30/09  
 Prep Date.....: 07/01/09 Analysis Date...: 07/03/09  
 Prep Batch #...: 9182192  
 Dilution Factor: 0.98 Method.....: SW846 8330

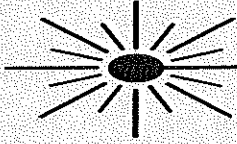
PARAMETER	RESULT	REPORTING LIMIT	UNITS	MDL
Nitroglycerin	ND	0.64	ug/L	0.15
PETN	ND	0.64	ug/L	0.23
Picric Acid	ND	0.98	ug/L	0.12
2-Amino-4,6-dinitrotoluene	ND	0.20	ug/L	0.098
4-Amino-2,6-dinitrotoluene	ND	0.098	ug/L	0.022
1,3-Dinitrobenzene	ND	0.098	ug/L	0.049
2,4-Dinitrotoluene	ND	0.098	ug/L	0.049
2,6-Dinitrotoluene	ND	0.098	ug/L	0.049
HMX	0.25 J	0.098	ug/L	0.026
Nitrobenzene	ND	0.098	ug/L	0.049
2-Nitrotoluene	ND	0.39	ug/L	0.071
3-Nitrotoluene	ND	0.39	ug/L	0.061
4-Nitrotoluene	ND	0.49	ug/L	0.071
RDX	4.1 J	0.098	ug/L	0.064
Tetryl	ND	0.098	ug/L	0.049
1,3,5-Trinitrobenzene	<del>0.038</del> J, B	0.098 U	ug/L	0.030
2,4,6-Trinitrotoluene	ND	0.098	ug/L	0.024

SURROGATE	PERCENT RECOVERY	RECOVERY LIMITS
3,4-Dinitrotoluene	105	(79 - 111)

NOTE(S):

- J Estimated result. Result is less than RL.
- B Method blank contamination. The associated method blank contains the target analyte at a reportable level.

*WS 8/6/09*



ECT.CON INC.

Environmental and Computer  
Technology Consultants

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## *Support Documentation*

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## *Laboratory Case Narrative*



## **DATA DELIVERABLES PACKAGE**

Michael Baker Jr., Inc.  
James D. Peyton  
5261 Fountain Drive, Suite A  
Crown Point, IN 46307

Client Project: Camp Bonneville Groundwater  
Client Project Number: 110002 Task 6200

TA Work Order#: PSF0891  
TA Project Manager: Estella Rieben

The total number of pages contained in this data package is:

187

July 31, 2009

TestAmerica - Portland  
9405 S. W. Nimbus Avenue  
Beaverton, Oregon 97008  
(503) 906-9200  
(503) 906-9210

## CASE NARRATIVE

**Client:** Michael Baker Jr., Inc.                      **Date Sampled:** 6/25/2009 to 6/26/2009  
**Project:** Camp Bonneville Groundwater              **Date Received:** 6/26/2009  
                  110002 Task 6200  
**Lab:** PSF0891

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**SAMPLE RECEIPT:** Samples were received intact, on ice, with chain of custody documentation. The sample temperature was measured at 0.2 and 2.0 °C upon receipt at the laboratory.

**HOLDING TIMES:** All samples were analyzed within holding times.

**PROBLEMS ENCOUNTERED:** No problems were encountered.

**QA/QC CRITERIA:** 1,2,3-Trichlorobenzene was detected at a concentration less than the MRL but greater than one-half of the MRL in the associated method blank (9070189-BLK1) for EPA 8260B samples PSF0891-02, PSF0891-03, and PSF0891-04. The concentration for the specified analyte in all client samples was ND.

**OBSERVATIONS:** EPA 8260B sample PSF0891-01 was prepared in batch 9070189, analyzed on 07/08/09, but was not reported due to suspected contamination caused by carry-over from a previously analyzed sample. The specified sample was re-prepared in batch 9070249, designated "RE1," analyzed on 07/09/09, and reported.

**SUBCONTRACTED:** 314.0\_9196313 (Perchlorate), 314.0\_9196314 (Perchlorate), and SW846 8330 analyses were performed at the TestAmerica facility in West Sacramento, CA. The data package for the specified analyses is included in this data deliverables package as Appendix A.

**Michael Baker Jr., Inc.**  
5261 Fountain Drive, Suite A  
Crown Point, IN 46307

Project Name: **Camp Bonneville Groundwater**  
Project Number: 110002 Task 6200  
Project Manager: James D. Peyton

Report Created:  
07/21/09 13:15

## ANALYTICAL REPORT FOR SAMPLES

Sample ID	Laboratory ID	Matrix	Date Sampled	Date Received
23LCMW01SW	PSF0891-01	Water	06/25/09 16:30	06/26/09 16:50
23LF4MW17SW	PSF0891-02	Water	06/25/09 18:00	06/26/09 16:50
23LF4MW18SW	PSF0891-03	Water	06/25/09 19:00	06/26/09 16:50
23LF4MW7B	PSF0891-04	Water	06/26/09 10:30	06/26/09 16:50
23LF4MW5A	PSF0891-05	Water	06/26/09 13:45	06/26/09 16:50
TB261	PSF0891-06	Water	06/26/09 00:00	06/26/09 16:50

TestAmerica Portland



Estella Rieben, Project Manager

*The results in this report apply to the samples analyzed in accordance with the chain of custody document. This analytical report shall not be reproduced except in full, without the written approval of the laboratory.*



# TestAmerica

THE LEADER IN ENVIRONMENTAL TESTING

July 28, 2009

**TestAmerica Project Number: G9F300242**

PO/Contract: PSF0891

Estella Rieben  
TestAmerica Portland  
Nimbus Corporate Center  
9405 SW Nimbus Ave  
Beaverton, OR 97008

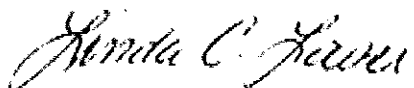
Dear Ms. Rieben,

This report contains the analytical results for the samples received under chain of custody by TestAmerica on June 30, 2009. These samples are associated with your Michael Baker Jr., Inc. project.

The test results in this report meet all NELAC requirements for parameters that accreditation is required or available. Any exceptions to NELAC requirements are noted in the case narrative. The case narrative is an integral part of this report.

If you have any questions, please feel free to call me at (916) 374-4362.

Sincerely,



Linda C. Laver  
Project Manager

## Case Narrative

### TestAmerica West Sacramento Project Number G9F300242

#### **WATER, 8330, Nitroaromatics & Nitramines**

Samples: 1, 2, 3, 4, 5

The laboratory control sample (LCS) associated with this extraction batch has percent recoveries for most analytes above the established control limits indicating a potential high bias in the data. Samples 1, 2, 3, and 4 do not have detected concentrations above the reporting limit for these analytes and there is no adverse impact upon the data. Sample 5 does have detected concentrations for HMX and RDX. This was not discovered until after the recommended hold time for extraction had expired. After consulting with Ms. Rieben, the sample was not re-extracted and the analytical results are included in this report.

The matrix spike/matrix spike duplicate (MS/MSD) associated with this extraction batch also has recoveries and/or precision is outside the established control limits for many analytes. This anomaly is most likely matrix related.

There are no other anomalies associated with this project.

## Sample Summary

### TestAmerica West Sacramento Project Number G9F300242

<u>WO#</u>	<u>Sample #</u>	<u>Client Sample ID</u>	<u>Sampling Date</u>	<u>Received Date</u>
LFV9A	1	PSF0891-01	6/25/2009 04:30 PM	6/30/2009 09:00 AM
LFV9G	2	PSF0891-02	6/25/2009 06:00 PM	6/30/2009 09:00 AM
LFV9J	3	PSF0891-03	6/25/2009 07:00 PM	6/30/2009 09:00 AM
LFV9K	4	PSF0891-04	6/26/2009 10:30 AM	6/30/2009 09:00 AM
LFV9M	5	PSF0891-05	6/26/2009 01:45 PM	6/30/2009 09:00 AM

#### Notes(s):

- The analytical results of the samples listed above are presented on the following pages.
- All calculations are performed before rounding to avoid round-off errors in calculated results.
- Results noted as "ND" were not detected at or above the stated limit.
- This report must not be reproduced, except in full, without the written approval of the laboratory.
- Results for the following parameters are never reported on a dry weight basis: color, corrosivity, density, flashpoint, ignitability, layers, odor, paint filter test, pH, porosity, pressure, reactivity, redox potential, specific gravity, spot tests, solids, solubility, temperature, viscosity, and weight.

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## *Chain of Custody*





TestAmerica Portland  
**Sample Receiving Checklist**

Work Order #: PSF0891 Date/Time Received: 4/20/09 1650  
 Client Name and Project: Michael DeKee - Camp Bonneville GW

Time Zone:  
 EDT/EST     CDT/CST     MDT/MST     PDT/PST     AK     OTHER

**Unpacking Checks:**

Cooler #(s): 1  
 Temperatures: 6 6  
 Digi #1  Digi #2  IR Gun  (  Plastic  Glass )

**Temperature out of Range:**

Not enough or No Ice  
 Ice Melted  
 W/in 4 Hrs of collection  
 Other: \_\_\_\_\_

Initials: PS

N.A Yes No

- 1. If ESI client, were temp blanks received? If no, document on NOD.
- 2. Cooler Seals intact? (N/A if hand delivered) if no, document on NOD.
- 3. Chain of Custody present? If no, document on NOD.
- 4. Bottles received intact? If no, document on NOD.
- 5. Sample is not multiphasic? If no, document on NOD.
- 6. Proper Container and preservatives used? If no, document on NOD.
- 7. pH of all samples checked and meet requirements? If no, document on NOD.
- 8. Cyanide samples checked for sulfides and meet requirements? If no, notify PM.
- 9. HF Dilution required?
- 10. Sufficient volume provided for all analysis? If no, document on NOD and consult PM before proceeding.
- 11. Did chain of custody agree with samples received? If no, document on NOD.
- 12. Is the "Sampled by" section of the COC completed?
- 13. Were VOA/Oil Syringe samples without headspace?
- 14. Were VOA vials preserved?  HCl  Sodium Thiosulfate  Ascorbic Acid
- 15. Did samples require preservation with sodium thiosulfate?
- 16. If yes to #14, was the residual chlorine test negative? If no, document on NOD.
- 17. Are dissolved/field filtered metals bottles sediment-free? If no, document on NOD.
- 18. Is sufficient volume provided for client requested MS/MSD or matrix duplicates? If no, document on NOD and contact PM before proceeding.
- 19. Are analyses with short holding times received in hold?
- 20. Was Standard Turn Around (TAT) requested?
- 21. Receipt date(s) < 48 hours past the collection date(s)? If no, notify PM.

TestAmerica Portland  
Sample Receiving Checklist

Work Order #: PSF0891

Login Checks:

Initials: BUE

N/A Yes No

22. Sufficient volume provided for all analysis? If no, document on NOD & contact PM.
23. Sufficient volume provided for client requested MS/MSD or matrix duplicates? If no, document on NOD and contact PM.
24. Did the chain of custody include "received by" and "relinquished by" signatures, dates and times?
25. Were special log in instructions read and followed?
26. Were tests logged checked against the COC?
27. Were rush notices printed and delivered?
28. Were short hold notices printed and delivered?
29. Were subcontract COCs printed?
30. Was HF dilution logged?

Labeling and Storage Checks:

Initials: PS

N/A Yes No

31. Were the subcontracted samples/containers put in Sx fridge?
32. Were sample bottles and COC double checked for dissolved/filtered metals?
33. Did the sample ID, Date, and Time from label match what was logged?
34. Were Foreign sample stickers affixed to each container and containers stored in foreign fridge?
35. Were HF stickers affixed to each container, and containers stored in Sx fridge?
36. Was an NOD for created for noted discrepancies and placed in folder?

Document any problems or discrepancies and the actions taken to resolve them on a Notice of Discrepancy form (NOD).

CLIENT TAL - Portland PM U LOG# 59265

LOT# (QUANTIMS ID) 69F30242 QUOTE# 83365 LOCATION WIB

DATE RECEIVED 6/30/09 TIME RECEIVED 0900 Initials CV Date 6/30/09

DELIVERED BY  FEDEX  CA OVERNIGHT  CLIENT  
 AIRBORNE  GOLDENSTATE  DHL  
 UPS  BAX GLOBAL  GO-GETTERS  
 TAL COURIER  VALLEY LOGISTICS  MORGAN HILL COURIER  
 OTHER

CUSTODY SEAL STATUS  INTACT  BROKEN  N/A

CUSTODY SEAL #(S) 204315, 204295

SHIPPING CONTAINER(S)  TAL  CLIENT  N/A

TEMPERATURE RECORD (IN °C) IR 4  5  OTHER \_\_\_\_\_

COC #(S) 10

TEMPERATURE BLANK Observed: 2 Corrected: 2  
006/30/09

SAMPLE TEMPERATURE

Observed: 2 1 4 Average: 3 Corrected Average: 3

COLLECTOR'S NAME: 2 3 4  Verified from COC  Not on COC 3

pH MEASURED  YES  ANOMALY  N/A

LABELED BY.....

LABELS CHECKED BY.....

PEER REVIEW  NA

SHORT HOLD TEST NOTIFICATION

SAMPLE RECEIVING

WETCHEM  N/A

VOA-ENCORES  N/A

METALS NOTIFIED OF FILTER/PRESERVE VIA VERBAL & EMAIL  N/A

COMPLETE SHIPMENT RECEIVED IN GOOD CONDITION WITH APPROPRIATE TEMPERATURES, CONTAINERS, PRESERVATIVES  N/A

CLOUSEAU  TEMPERATURE EXCEEDED (2 °C - 6 °C)\*1  N/A

WET ICE  BLUE ICE  GEL PACK  NO COOLING AGENTS USED  PM NOTIFIED

Notes: All RT req'd @ i's

\*1 Acceptable temperature range for State of Wisconsin samples is ≤4°C

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# *Worksheets*

**HOLDING TIMES**

SAMPLE DATE	SAMPLE ID	Portland LAB ID	Sacramento Lab ID	Matrix	Prep	Anal
06/25/09	23LCMW01SW	PSF0891-01	LFV9A	Aq	07/01/09	7/3/09 1X
06/25/09	23LF4MW17SW	PSF0891-02	LFV9G	Aq	07/01/09	7/3/09 1X
06/25/09	23LF4MW18SW	PSF0891-03	LFV9J	Aq	07/01/09	7/3/09 1X
06/26/09	23LF4MW7B	PSF0891-04	LFV9K	Aq	07/01/09	7/3/09 1X
06/26/09	23LF4MW5A	PSF0891-05	LFV9M	Aq	07/01/09	7/3/09 1X

7 days from collection to extraction  
 Sample Date 6/25/2009 Extract By 7/2/2009

7 days from collection to extraction  
 Sample Date 6/26/2009 Extract By 7/3/2009

40 days from extraction to analysis  
 Sample Date 7/1/2009 Extract By 8/10/2009

**TARGET COMPOUNDS AND QUANTITATION LIMITS**

Was a target compound list provided by the client? No  
 Did Sample Form 1s match the target compound list NA  
 Were required quantitation limits provided by the client? No  
 Did all compounds meet the required quantitation limits? NA

**SYSTEM MONITORING COMPOUNDS**

SAMPLE	SURROGATE	COLUMN 1	DF	ACTION
All IN				

Were surrogate RTs within windows established by the ICAL? Yes  
 Were there any transcription errors between the raw data and Form 2? No  
 Were laboratory acceptance limits used as the basis for validation? 30-150  
 Did the laboratory provide CLP Form II or equivalent? Yes

MW5A

SURR	AMOUNT FOUND	AMOUNT SPIKED	% R	F1
34-DNT A	2.588	2.5	103.5	105
34-DNT C	2.81	2.5	112.4	

**MS/MSD RECOVERY and RELATIVE PERCENT DIFFERENCE**

23LF4MW7B						
COMPOUND	MS %R	MSD %R	RPD	NATIVE	SPIKE	ACTION
picric acid	ok	ok	35	ND		note
2-a-46-dnt	125	129	ok	ND		note
4-a-26-dnt	116	120	ok	ND		note
13dmben	125	130	ok	ND		note
24-dnt	120	124	ok	ND		note
26-dnt	122	125	ok	ND		note
hmx	118	120	ok	ND		note
nbz	121	124	ok	ND		note
3-nt	ok	115	ok	ND		note
4-nt	ok	116	ok	ND		note
rdx	ok	129	ok	ND		note
tetryl	ok	106	ok	ND		note
135-tnb	ok	124	ok	ND		note
246-tnt	113	117	ok	ND		note

Were there any transcription errors between the raw data and Form 3? No  
 Were laboratory acceptance limits used as the basis for validation? Yes  
 Did the laboratory provide CLP Form III or equivalent? Yes  
 Were chromatograms and quan reports present for all MS/MSDs? Yes

picric acid	AMOUNT FOUND	AMOUNT NATIVE	AMOUNT SPIKED	%R	Form 3
MS	3.41	0	4.91	69.5	70
MSD	4.85	0	4.92	98.6	99
RPD	-34.9			-34.7	35.0

Recovery = ((Amount Found-Amount Native)/Amount Spiked)\*100  
 RPD = ((MS-MSD)/((MS+MSD)/2))\*100

**LABORATORY CONTROL SAMPLES**

SAMPLE	COMPOUND	LCS %R	LCSD %R	RPD	ACTION
LFW791	12 of 17 HI	OK	NA	NA	J

J HMX and RDX in 23LF4MW5A

Were laboratory acceptance limits used as the basis for validation? Yes  
 Did the laboratory provide CLP Form III or equivalent? Yes  
 Were chromatograms and quan reports present for all LCS/LCDs? Yes

LFW791	RDX	AMOUNT FOUND	AMOUNT SPIKED	% R	FORM 3
12 of 17 HI					
LCS		1.27	1	127.0	127

%R = (Amount Found/Amount Spiked)\*100

**BLANKS**

BLANK	COMPOUND	RESULT	5X OR 10X	ACTION LEVEL	ACTION
LFW791	135-tmbz	0.039	5	0.195	U at RL

NOTE Equipment/Field blanks are not qualified on the basis of laboratory method blank contamination or contamination in other field quality control blanks.

**CALIBRATION**

**Initial**

Are chromatograms and quan reports present for all ICAL standards? Yes  
 Are CLP For VIs or equivalent present and complete? Yes  
 Any transcription or calculation errors? No  
 What linearity criteria was used? 20% RSD or 0.99  
 Were RT window documented? Yes

**Continuing**

Are chromatograms and quan reports present for all CCV standards? Yes  
 Are CLP Form VIIs or equivalent present and complete? Yes  
 Any transcription or calculation errors? No  
 What %D criteria was used? 25% D  
 Have all samples and standards been listed on an Analytical Sequence? Yes  
 Was a proper analytical sequence followed? Yes

**INITIAL CALIBRATION - A**

Date	6/10/2009	Rptd Avg CF	77.89163	Rptd %RSD	5.156
Instrument	A	Calcd Avg CF	77.892	Calcd %RSD	5.156
Compound	RDX				
Reported CF	80.13	CF1	81.80	STD DEV	4
Calculated CF	80.42	CF2	75.90		
		CF3	77.15		
Response	16084	CF4	80.72		
Conc	200	CF5	78.45		
		CF6	80.13		
	CF = (H/C)	CF7	69.168	%RSD = (Std Dev/Avg CF)*100	
		CF8	79.815		

**CONTINUING CALIBRATION A**

Date	7/3/2009		
Time	0045		
Instrument	A		
Compound	HMX		
Reported Conc	104.8	Rptd %D	5
Calculated Conc	103.835	Calcd %D	-3.83
Response	12564	ICAL CF	100
CF	121	CCV CF	103.8
	CF = (H/C)	%Difference = ((ICAL - CCV)/ICAL)*100	

**INITIAL CALIBRATION - C**

Date	6/9/2009	Rptd Avg CF	42.55812	Rptd %RSD	4.852
Instrument	C	Calcd Avg CF	42.558	Calcd %RSD	4.852
Compound	RDX				
Reported CF	42.585	CF1	45.40	STD DEV	2
Calculated CF	42.78	CF2	43.70		
		CF3	42.75		
Response	8556	CF4	43.30		
Conc	200	CF5	42.16		
		CF6	42.59		
	CF = (H/C)	CF7	38.118	%RSD = (Std Dev/Avg CF)*100	
		CF8	42.452		

**CONTINUING CALIBRATION C**

Date	7/11/2009		
Time	0554		
Instrument	C		
Compound	HMX		
Reported Conc	184.3	Rptd %D	8
Calculated Conc	184.280	Calcd %D	7.86
Response	7594	ICAL CF	200
CF	41.20913	CCV CF	184.3
	CF = (H/C)	%Difference = ((ICAL - CCV)/ICAL)*100	

**FIELD DUPLICATES**

COMPOUND	SAMPLE	QUALIFIER	DUPLICATE	QUALIFIER	RPD
PCA					#DIV/0!

NOTES Samples are not qualified on this basis.

**COMPOUND IDENTIFICATION AND QUANTITATION**

Has a F10 been completed for every sample containing positive results?	Yes
Was RT data presented on the form?	NA
Are RTs within the established windows?	Yes
Any transcription or calculation errors?	No
Any false positives, negative peaks, shouldering, etc.?	No
Was GCMS confirmation needed for results > 10 µg/ml?	NA
Were percent differences or relative percent differences calculated?	NA
Are percent differences/RPDs greater than 25%?	NA
Are there any transcription errors?	No
Are Form 1s present for all field and quality control samples?	No
Are chromatograms and quan reports present for all samples?	Yes
Are RLs adjusted to reflect sample dilutions, percent solids, etc.?	Yes
For soils, any percent solids <50%?	No
For soils, any percent solids <10%?	No

23LF4MW5A RDX 0.41 µg/L

	A	C
response	16156	8165
cf	77.89163	42.55812
final vol ml	20	20
initial L	1.01672	1.01673
df	1	1
calculated	4.0801	3.7740

**PERCENT DIFFERENCE BETWEEN COLUMNS**

SAMPLE	COMPOUND	A	C	RPD	QUALIFIER
23LF4MW5A	HMX	0.246	0.2052	18.1	none
	RDX	4.08	3.774	7.8	none

MATRIX SPIKE SAMPLE EVALUATION REPORT

HPLC

Client Lot #...: G9F300242      Work Order #...: Lfv9K1AF-MS      Matrix.....: WATER  
 MS Lot-Sample #: G9F300242-004      Lfv9K1AG-MSD  
 Date Sampled...: 06/26/09      Date Received...: 06/30/09  
 Prep Date.....: 07/01/09      Analysis Date...: 07/03/09  
 Prep Batch #...: 9182192  
 Dilution Factor: 0.98

23LF4MW7B (ND)

PARAMETER	PERCENT RECOVERY	RECOVERY LIMITS	RPD	RPD LIMITS	METHOD
Nitroglycerin	106	(85 - 115)			SW846 8330
	108	(85 - 115)	2.0	(0-15)	SW846 8330
PETN	110	(84 - 117)			SW846 8330
	111	(84 - 117)	0.90	(0-15)	SW846 8330
Picric Acid	70	(21 - 118)			SW846 8330
	99 p	(21 - 118)	35	(0-20)	SW846 8330
2-Amino-4,6-dinitrotoluene	125 a	(77 - 123)			SW846 8330
	129 a	(77 - 123)	3.4	(0-27)	SW846 8330
4-Amino-2,6-dinitrotoluene	116 a	(68 - 113)			SW846 8330
	120 a	(68 - 113)	3.0	(0-30)	SW846 8330
1,3-Dinitrobenzene	125 a	(72 - 123)			SW846 8330
	130 a	(72 - 123)	4.1	(0-29)	SW846 8330
2,4-Dinitrotoluene	120 a	(70 - 119)			SW846 8330
	124 a	(70 - 119)	2.9	(0-30)	SW846 8330
2,6-Dinitrotoluene	122 a	(71 - 119)			SW846 8330
	125 a	(71 - 119)	2.9	(0-29)	SW846 8330
HMX	118 a	(67 - 115)			SW846 8330
	120 a	(67 - 115)	1.4	(0-32)	SW846 8330
Nitrobenzene	121 a	(69 - 119)			SW846 8330
	124 a	(69 - 119)	2.5	(0-31)	SW846 8330
2-Nitrotoluene	117	(64 - 120)			SW846 8330
	113	(64 - 120)	2.7	(0-36)	SW846 8330
3-Nitrotoluene	113	(67 - 114)			SW846 8330
	115 a	(67 - 114)	2.0	(0-31)	SW846 8330
4-Nitrotoluene	114	(67 - 115)			SW846 8330
	116 a	(67 - 115)	2.3	(0-32)	SW846 8330
RDX	121	(68 - 122)			SW846 8330
	129 a	(68 - 122)	6.3	(0-32)	SW846 8330
Tetryl	102	(66 - 105)			SW846 8330
	106 a	(66 - 105)	4.2	(0-26)	SW846 8330
1,3,5-Trinitrobenzene	119	(74 - 120)			SW846 8330
	124 a	(74 - 120)	4.3	(0-29)	SW846 8330
2,4,6-Trinitrotoluene	113 a	(69 - 111)			SW846 8330
	117 a	(69 - 111)	3.4	(0-28)	SW846 8330

(Continued on next page)



MATRIX SPIKE SAMPLE EVALUATION REPORT

HPLC

Client Lot #...: G9F300242      Work Order #...: LFV9K1AF-MS      Matrix.....: WATER  
MS Lot-Sample #: G9F300242-004      LFV9K1AG-MSD

<u>SURROGATE</u>	<u>PERCENT</u>	<u>RECOVERY</u>
	<u>RECOVERY</u>	<u>LIMITS</u>
3,4-Dinitrotoluene	104	(79 - 111)
	107	(79 - 111)

**NOTE (S) :**

Calculations are performed before rounding to avoid round-off errors in calculated results

Bold print denotes control parameters

p Relative percent difference (RPD) is outside stated control limits

a Spiked analyte recovery is outside stated control limits.

MATRIX SPIKE SAMPLE DATA REPORT

HPLC

Client Lot #...: G9F300242      Work Order #...: LFV9K1AF-MS      Matrix.....: WATER  
 MS Lot-Sample #: G9F300242-004      LFV9K1AG-MSD  
 Date Sampled...: 06/26/09      Date Received...: 06/30/09  
 Prep Date.....: 07/01/09      Analysis Date...: 07/03/09  
 Prep Batch #...: 9182192  
 Dilution Factor: 0.98

PARAMETER	SAMPLE AMOUNT	SPIKE AMT	MEASRD AMOUNT	UNITS	PERCNT RECVRY	RPD	METHOD
Nitroglycerin	ND	4.91	5.19	ug/L	106		SW846 8330
	ND	4.92	5.29	ug/L	108	2.0	SW846 8330
PETN	ND	4.91	5.41	ug/L	110		SW846 8330
	ND	4.92	5.46	ug/L	111	0.90	SW846 8330
Picric Acid	ND	4.91	3.41	ug/L	70		SW846 8330
	ND	4.92	4.85	ug/L	99 p	35	SW846 8330
2-Amino-4,6-dinitrotoluene	ND	0.982	1.22	ug/L	125 a		SW846 8330
	ND	0.983	1.26	ug/L	129 a	3.4	SW846 8330
4-Amino-2,6-dinitrotoluene	ND	0.982	1.14	ug/L	116 a		SW846 8330
	ND	0.983	1.18	ug/L	120 a	3.0	SW846 8330
1,3-Dinitrobenzene	ND	0.982	1.23	ug/L	125 a		SW846 8330
	ND	0.983	1.28	ug/L	130 a	4.1	SW846 8330
2,4-Dinitrotoluene	ND	0.982	1.18	ug/L	120 a		SW846 8330
	ND	0.983	1.22	ug/L	124 a	2.9	SW846 8330
2,6-Dinitrotoluene	ND	0.982	1.19	ug/L	122 a		SW846 8330
	ND	0.983	1.23	ug/L	125 a	2.9	SW846 8330
HMX	ND	0.982	1.16	ug/L	118 a		SW846 8330
	ND	0.983	1.18	ug/L	120 a	1.4	SW846 8330
Nitrobenzene	ND	0.982	1.19	ug/L	121 a		SW846 8330
	ND	0.983	1.22	ug/L	124 a	2.5	SW846 8330
2-Nitrotoluene	ND	0.982	1.14	ug/L	117		SW846 8330
	ND	0.983	1.11	ug/L	113	2.7	SW846 8330
3-Nitrotoluene	ND	0.982	1.11	ug/L	113		SW846 8330
	ND	0.983	1.14	ug/L	115 a	2.0	SW846 8330
4-Nitrotoluene	ND	0.982	1.12	ug/L	114		SW846 8330
	ND	0.983	1.14	ug/L	116 a	2.3	SW846 8330
RDX	ND	0.982	1.19	ug/L	121		SW846 8330
	ND	0.983	1.27	ug/L	129 a	6.3	SW846 8330
Tetryl	ND	0.982	0.998	ug/L	102		SW846 8330
	ND	0.983	1.04	ug/L	106 a	4.2	SW846 8330
1,3,5-Trinitrobenzene	0.035	0.982	1.20	ug/L	119		SW846 8330
	0.035	0.983	1.26	ug/L	124 a	4.3	SW846 8330
2,4,6-Trinitrotoluene	ND	0.982	1.11	ug/L	113 a		SW846 8330
	ND	0.983	1.15	ug/L	117 a	3.4	SW846 8330

(Continued on next page)

MATRIX SPIKE SAMPLE DATA REPORT

HPLC

Client Lot #...: G9F300242      Work Order #...: LFV9K1AF-MS      Matrix.....: WATER  
MS Lot-Sample #: G9F300242-004      LFV9K1AG-MSD

<u>SURROGATE</u>	<u>PERCENT RECOVERY</u>	<u>RECOVERY LIMITS</u>
3,4-Dinitrotoluene	104	(79 - 111)
	107	(79 - 111)

**NOTE (S) :**

Calculations are performed before rounding to avoid round-off errors in calculated results.

Bold print denotes control parameters

**p** Relative percent difference (RPD) is outside stated control limits.

**a** Spiked analyte recovery is outside stated control limits

MS

Chromatography Summary

Injection Date: 7/3/2009 6:39 Operator: fhk  
 DataFile: LC10 I07022009.BVA-000019.D Vial Num: 25  
 Instrument ID: LC10

Method 8330 Target Analyte Results

Sample: **LFV9K1AF 9182192 G9F300242-4 S 1X**

Method File: LC10 I07022009.BV8330AB.M  
 Start Cal Date: 6/10/2009 15:33 End Cal Date: 6/11/2009 13:47

Matrix: WATER SubList: WATER.sub SpikeList: WATER.spk  
 Samp. Info: LFV9K1AF 9182192 G9F300242-4 S 1X;3;  
 Misc. Info: LCS, 1018 28;;20;1;WATER sub,WATER spk,1,1,LFV9K1AF

Dilution Factor	Extract Volume	Sample Volume	Sample Weight
1X	20 mL	1018.28 mL	0 g

Signal 1 UV 250-265								Signal 2 UV 358-205								
Compound Name	RT	Response	Conc (ug/L)	Spike Level	%R	Result	Flag	RT	Response	Conc (ug/L)	Spike Level	%R	Result	Flag	Limits(%)	Flag
3,4-Dinitrotoluene	18.25	6156	2.5630	< 2.45512	104%	Acceptable		18.25	11386	2.6320	2.45512	107%	Acceptable	M	(79-111)	
HMX	5.30	7138	1.1600	< 0.982048	118%	Fails					0.982048	0%	Fails		(65-115)	45
RDX	7.88	4720	1.1900	< 0.982048	121%	Acceptable					0.982048	0%	Fails		(68-122)	45
Picric ACID	9.09	14543	3.4510	< 4.910241	70%	Acceptable		9.09	21152	3.4140	< 4.910241	70%	Acceptable		(21-118)	
1,3,5-Trinitrobenzene	10.24	9036	1.2020	< 0.982048	122%	Fails					0.982048	0%	Fails		(74-120)	45
1,3-Dinitrobenzene	13.31	8723	1.2260	< 0.982048	125%	Fails					0.982048	0%	Fails		(72-123)	45
TETRYL	14.37	4713	0.9976	< 0.982048	102%	Acceptable					0.982048	0%	Fails		(66-105)	45
Nitrobenzene	15.23	3678	1.1910	< 0.982048	121%	Fails					0.982048	0%	Fails		(69-119)	45
2,4,6-Trinitrotoluene	16.60	4967	1.1120	< 0.982048	113%	Fails					0.982048	0%	Fails		(69-111)	45
4-AM-2,6-DNT	17.61	3628	1.1440	< 0.982048	116%	Fails					0.982048	0%	Fails		(68-113)	45
2-AM-4,6-DNT	18.67	4580	1.2230	< 0.982048	125%	Fails					0.982048	0%	Fails		(77-123)	45
2,6-Dinitrotoluene	20.41	3237	1.1940	< 0.982048	122%	Fails					0.982048	0%	Fails		(71-119)	45
2,4-Dinitrotoluene	21.13	5152	1.1820	< 0.982048	120%	Fails					0.982048	0%	Fails		(70-119)	45
2-Nitrotoluene	24.86	2095	1.1450	< 0.982048	117%	Acceptable					0.982048	0%	Fails		(64-120)	45
4-Nitrotoluene	26.68	2486	1.1150	< 0.982048	114%	Acceptable					0.982048	0%	Fails		(67-115)	45
3-Nitrotoluene	28.70	2444	1.1130	< 0.982048	113%	Acceptable					0.982048	0%	Fails		(67-114)	45
Nitroglycerin				4.910241	0%	Fails		15.86	16078	5.1910	< 4.910241	106%	Acceptable		(85-115)	45
PETN				4.910241	0%	Fails		30.30	8828	5.4100	< 4.910241	110%	Acceptable		(84-117)	45
3,5-Dinitroaniline	14.05	5434	1.1860	< 0.982048	121%	Acceptable					0.982048	0%	Fails		(40-140)	45

Surrogates:	Spiked	Recovered	% Rec	Spiked	Recovered	% Rec	Limits
3,4-Dinitrotoluene	2.4551	2.5630	104	2.4551	2.6320	107	(79-111)

Notes: M = Manually Integrated Signals Differ by More Than 40%  
 D = Operator Disabled Result Signals Differ by More Than 50%  
 O = Over Calibration Range

RE-INTEGRATION CODES

- 1 Poor Peak Shape
  - 2 Poor Peak Resolution
  - 3 Peak Not Integrated
  - 4 Sample Matrix Interference
  - 5 Column Bleed
  - 6 Instrument Noise
  - 7 Baseline Correction
  - 8 Other (reason must be stated)
- ALL RE-INTEGRATIONS MUST BE INITIALED, DATED AND CODED

*fhk 7/7/09*

MSD

A-000020.D

Chromatography Summary

Injection Date: 7/3/2009 7:30 Operator: fhk  
 DataFile: LC10 I07022009 BVA-000020.D Vial Num: 26  
 Instrument ID: LC10

Method 8330 Target Analyte Results

Sample: **LFV9KIAG 9182192 G9F300242-4 D  
 1X**

Method File: LC10 I07022009 B8330AB.M  
 Start Cal Date: 6/10/2009 15:33 End Cal Date: 6/11/2009 13:47

Matrix: WATER SubList: WATER.sub SpikeList: WATER.spk  
 Samp. Info: LFV9KIAG 9182192 G9F300242-4 D 1X,3  
 Misc. Info: LCS,,1017 3,,20,1;WATER.sub,WATER.spk,1,1,LFV9KIAG

Dilution Factor	Extract Volume	Sample Volume	Sample Weight
1X	20 mL	1017.3 mL	0 g

Signal 1 UV 250-265								Signal 2 UV 358-205								
Compound Name	RT	Response	Conc (ug/L)	Spike Level	%R	Result	Flag	RT	Response	Conc (ug/L)	Spike Level	%R	Result	Flag	Limits(%)	Flag
3,4-Dinitrotoluene	18.25	6293	2.6230	< 2.457486	107%	Acceptable		18.25	10744	2.4430	2.457486	99%	Acceptable		(79-111)	
HMX	5.31	7227	1.1760	< 0.982994	120%	Fails					0.982994	0%	Fails		(65-115)	45
RDX	7.89	5021	1.2670	< 0.982994	129%	Fails					0.982994	0%	Fails		(68-122)	45
Picric ACID	9.08	20473	4.8620	< 4.914971	99%	Acceptable		9.08	29993	4.8460	< 4.914971	99%	Acceptable		(21-118)	
1,3,5-Trinitrobenzene	10.24	9422	1.2550	< 0.982994	128%	Fails					0.982994	0%	Fails		(74-120)	45
1,3-Dinitrobenzene	13.31	9074	1.2770	< 0.982994	130%	Fails					0.982994	0%	Fails		(72-123)	45
TETRYL	14.37	4912	1.0410	< 0.982994	106%	Fails					0.982994	0%	Fails		(66-105)	45
Nitrobenzene	15.25	3769	1.2210	< 0.982994	124%	Fails					0.982994	0%	Fails		(69-119)	45
2,4,6-Trinitrotoluene	16.60	5134	1.1510	< 0.982994	117%	Fails					0.982994	0%	Fails		(69-111)	45
4-AM-2,6-DNT	17.61	3736	1.1790	< 0.982994	120%	Fails					0.982994	0%	Fails		(68-113)	45
2-AM-4,6-DNT	18.68	4734	1.2650	< 0.982994	129%	Fails					0.982994	0%	Fails		(77-123)	45
2,6-Dinitrotoluene	20.40	3327	1.2290	< 0.982994	125%	Fails					0.982994	0%	Fails		(71-119)	45
2,4-Dinitrotoluene	21.13	5301	1.2170	< 0.982994	124%	Fails					0.982994	0%	Fails		(70-119)	45
2-Nitrotoluene	24.85	2037	1.1140	< 0.982994	113%	Acceptable					0.982994	0%	Fails		(64-120)	45
4-Nitrotoluene	26.67	2543	1.1410	< 0.982994	116%	Fails					0.982994	0%	Fails		(67-115)	45
3-Nitrotoluene	28.69	2491	1.1350	< 0.982994	115%	Fails					0.982994	0%	Fails		(67-114)	45
Nitroglycerin				4.914971	0%	Fails		15.86	16383	5.2940	< 4.914971	108%	Acceptable		(85-115)	45
PETN				4.914971	0%	Fails		30.29	8899	5.4590	< 4.914971	111%	Acceptable		(84-117)	45
3,5-Dinitroaniline	14.06	5660	1.2370	< 0.982994	126%	Acceptable					0.982994	0%	Fails		(40-140)	45

Surrogates:	Spiked	Recovered	%Rec	Spiked	Recovered	%Rec	Limits
3,4-Dinitrotoluene	2.4575	2.6230	107	2.4575	2.4430	99	(79-111)

Notes: M = Manually Integrated Signals Differ by More Than 40%  
 D = Operator Disabled Result Signals Differ by More Than 50%  
 O = Over Calibration Range

ed)

See reintegration.

*fhk*  
 7/2/09

LABORATORY CONTROL SAMPLE EVALUATION REPORT

HPLC

Client Lot #...: G9F300242      Work Order #...: LFW791AC      Matrix.....: WATER  
 LCS Lot-Sample#: G9G010000-192  
 Prep Date.....: 07/01/09      Analysis Date...: 07/02/09  
 Prep Batch #...: 9182192  
 Dilution Factor: 1

<u>PARAMETER</u>	<u>PERCENT RECOVERY</u>	<u>RECOVERY LIMITS</u>	<u>METHOD</u>
Nitroglycerin	107	(85 - 115)	SW846 8330
PETN	110	(84 - 117)	SW846 8330
Picric Acid	88	(21 - 118)	SW846 8330
2-Amino-4,6-dinitrotoluene	127 a	(77 - 123)	SW846 8330
4-Amino-2,6-dinitrotoluene	119 a	(68 - 113)	SW846 8330
1,3-Dinitrobenzene	128 a	(72 - 123)	SW846 8330
2,4-Dinitrotoluene	123 a	(70 - 119)	SW846 8330
2,6-Dinitrotoluene	125 a	(71 - 119)	SW846 8330
HMX	120 a	(67 - 115)	SW846 8330
Nitrobenzene	124 a	(69 - 119)	SW846 8330
2-Nitrotoluene	115	(64 - 120)	SW846 8330
3-Nitrotoluene	118 a	(67 - 114)	SW846 8330
4-Nitrotoluene	117 a	(67 - 115)	SW846 8330
RDX	127 a	(68 - 122)	SW846 8330
Tetryl	105	(66 - 105)	SW846 8330
1,3,5-Trinitrobenzene	127 a	(74 - 120)	SW846 8330
2,4,6-Trinitrotoluene	116 a	(69 - 111)	SW846 8330
<u>SURROGATE</u>		<u>PERCENT RECOVERY</u>	<u>RECOVERY LIMITS</u>
3,4-Dinitrotoluene		106	(79 - 111)

**NOTE(S) :**

Calculations are performed before rounding to avoid round-off errors in calculated results.

Bold print denotes control parameters

a Spiked analyte recovery is outside stated control limits.

LABORATORY CONTROL SAMPLE DATA REPORT

HPLC

Client Lot #...: G9F300242      Work Order #...: LFW791AC      Matrix.....: WATER  
 LCS Lot-Sample#: G9G010000-192  
 Prep Date.....: 07/01/09      Analysis Date...: 07/02/09  
 Prep Batch #...: 9182192  
 Dilution Factor: 1

<u>PARAMETER</u>	<u>SPIKE</u> <u>AMOUNT</u>	<u>MEASURED</u> <u>AMOUNT</u>	<u>UNITS</u>	<u>PERCENT</u> <u>RECOVERY</u>	<u>METHOD</u>
Nitroglycerin	5.00	5.35	ug/L	107	SW846 8330
PETN	5.00	5.52	ug/L	110	SW846 8330
Picric Acid	5.00	4.42	ug/L	88	SW846 8330
2-Amino-4,6-dinitrotoluene	1.00	1.27 a	ug/L	127	SW846 8330
4-Amino-2,6-dinitrotoluene	1.00	1.19 a	ug/L	119	SW846 8330
1,3-Dinitrobenzene	1.00	1.28 a	ug/L	128	SW846 8330
2,4-Dinitrotoluene	1.00	1.23 a	ug/L	123	SW846 8330
2,6-Dinitrotoluene	1.00	1.25 a	ug/L	125	SW846 8330
HMX	1.00	1.20 a	ug/L	120	SW846 8330
Nitrobenzene	1.00	1.24 a	ug/L	124	SW846 8330
2-Nitrotoluene	1.00	1.15	ug/L	115	SW846 8330
3-Nitrotoluene	1.00	1.18 a	ug/L	118	SW846 8330
4-Nitrotoluene	1.00	1.17 a	ug/L	117	SW846 8330
RDX	1.00	1.27 a	ug/L	127	SW846 8330
Tetryl	1.00	1.05	ug/L	105	SW846 8330
1,3,5-Trinitrobenzene	1.00	1.27 a	ug/L	127	SW846 8330
2,4,6-Trinitrotoluene	1.00	1.16 a	ug/L	116	SW846 8330

<u>SURROGATE</u>	<u>PERCENT</u> <u>RECOVERY</u>	<u>RECOVERY</u> <u>LIMITS</u>
3,4-Dinitrotoluene	106	(79 - 111)

**NOTE(S) :**

Calculations are performed before rounding to avoid round-off errors in calculated results

Bold print denotes control parameters

a Spiked analyte recovery is outside stated control limits.

METHOD BLANK REPORT

HPLC

Client Lot #...: G9F300242      Work Order #...: LEW791AA      Matrix.....: WATER  
 MB Lot-Sample #: G9C010000-192  
 Analysis Date...: 07/02/09      Prep Date.....: 07/01/09  
 Dilution Factor: 1      Prep Batch #...: 9182192

PARAMETER	RESULT	REPORTING		
		LIMIT	UNITS	METHOD
Nitroglycerin	ND	0.65	ug/L	SW846 8330
PETN	ND	0.65	ug/L	SW846 8330
Picric Acid	ND	1.0	ug/L	SW846 8330
2-Amino-4,6-dinitrotoluene	ND	0.20	ug/L	SW846 8330
4-Amino-2,6-dinitrotoluene	ND	0.10	ug/L	SW846 8330
1,3-Dinitrobenzene	ND	0.10	ug/L	SW846 8330
2,4-Dinitrotoluene	ND	0.10	ug/L	SW846 8330
2,6-Dinitrotoluene	ND	0.10	ug/L	SW846 8330
HMX	ND	0.10	ug/L	SW846 8330
Nitrobenzene	ND	0.10	ug/L	SW846 8330
2-Nitrotoluene	ND	0.40	ug/L	SW846 8330
3-Nitrotoluene	ND	0.40	ug/L	SW846 8330
4-Nitrotoluene	ND	0.50	ug/L	SW846 8330
RDX	ND	0.10	ug/L	SW846 8330
Tetryl	ND	0.10	ug/L	SW846 8330
1,3,5-Trinitrobenzene	0.039 J	0.10	ug/L	SW846 8330
2,4,6-Trinitrotoluene	ND	0.10	ug/L	SW846 8330

SURROGATE	PERCENT	RECOVERY
	RECOVERY	LIMITS
3,4-Dinitrotoluene	106	(79 - 111)

NOTE(S):

Calculations are performed before rounding to avoid round-off errors in calculated results.

J Estimated result. Result is less than RL.



# QC DATA ASSOCIATION SUMMARY

G9F300242

Sample Preparation and Analysis Control Numbers

<u>SAMPLE#</u>	<u>MATRIX</u>	<u>ANALYTICAL METHOD</u>	<u>LEACH BATCH #</u>	<u>PREP BATCH #</u>	<u>MS RUN#</u>
001	WATER	SW846 8330		9182192	9182194
002	WATER	SW846 8330		9182192	9182194
003	WATER	SW846 8330		9182192	9182194
004	WATER	SW846 8330		9182192	9182194
005	WATER	SW846 8330		9182192	9182194

TestAmerica West Sacramento  
ESC-Extraction Master Sheet

BI

Holding Time Due: 7/1/09 7/02/09  
BATCH #: 9182192  
Test #: 8330-L

Project Due: 7/10/09 7/13/09  
Initiated By: TP  
Extn Comp'd By: TP

Date: 7/01/09  
Date: 7/02/09

QC Code	Lab ID	Sample Size	Initial Mass	Final Volume	Final Mass	Difference in Mass
C	MB	See weight sheet	TP7/02/09			
	LCS					
	G9F260326-01					
	G9F260326-02					
	G9F260326-03					
7/01/09	G9F260270189-01					
	G9F270189-02					
	G9F270189-03					
	G9F270189-04					
	G9F270189-05					
	G9F300242-01					
	G9F300242-02					
	G9F300242-03					
S	G9F300242-04					
	G9F300242-04FS					
D	G9F300242-04NSD					
	G9F300242-05					
	TP 7/02/09					

SOP No: W5-LC-6509  
EXTRACTION COMMENTS:

Standard Information

QC Codes	Volume	STD ID	Exp. Date	STD Name/Conc.	ppm(ppb)
All	50uL	09GCSV0172	12/10/09	8330 SUCR	50ug/ml 2.5
GSD	20uL	09GCSV0037	7/17/09	3.5 DNA	50ug/ml 1.0
GSD	100uL	09GCSV0149	11/1/09	NG/PETN	50ug/ml 5.0
GSD	100uL	09GCSV0039	7/17/09	Picric ACID	50ug/ml 5.0

Spiked By / Date: TP 7/01/09  
Witnessed By / Date: ND 7-1-09

QC Codes: B = MB, C = LCS, L = LCSD, S = MS, D = SD, FB = Filter Blank

RQC058

TestAmerica Laboratories, Inc.  
EXTRACTION BENCH WORKSHEET

Run Date: 7/02/09  
Time: 11:49:39

LEV LEV LEV LEV  
1 2 1 2  
Y Y Y Y  
Y Y Y Y  
Y Y Y Y  
Y Y Y Y

Blank  
 Check  
 MS/MSD

Weights/Volumes  
 Spike & Surrogate Worksheet  
 Vial contains correct volume  
 Labels, greenbars, worksheets  
 computer batch: correct & all match  
 Anomalies to Extraction Method

Expanded Deliverable  
 COC Completed  
 Bench Sheet Copied  
 Package Submitted to Analytical Group  
 Bench Sheet Copied per COC

Extractionist: 002448 Tuan Q. Phan

Concentrationist: 002448 Tuan Q. Phan

\*\*\*\*\*  
 \* QC BATCH: 9182192 \*  
 \* PREP DATE: 7/01/09 9:00  
 \* COMP DATE: 7/02/09 10:40  
 \*\*\*\*\*

Reviewer/Date: PHANT / 7/02/09

Nitroaromatics & Nitramines: Explosives (8330)  
SOLID PHASE EXTRACTION (NOMINAL)

EXTR EXPR	ANL DUE	LOT# WORK ORDER	MSRUN#/ TEST FLGS	EXT MTH	MATRIX	INIT/ WT/VOL	PH'S ADJ1	EXTRACTION VOL	SOLVENTS EXCHANGE	VOL	SPIKE STANDARD/ SURROGATE ID		
7/01/09	7/10/09	G9F260326-001 LFOAK-1-AC		B7	A0 WATER	997.78mL 20.00mL	NA	NA	HOAC/ACN	4.5	HOAC/ACN	5.0	S0UL-09GCSV0172
7/01/09	7/10/09	G9F260326-002 LFOAL-1-AC		B7	A0 WATER	991.12mL 20.00mL	NA	NA	HOAC/ACN	4.5	HOAC/ACN	5.0	S0UL-09GCSV0172
7/01/09	7/10/09	G9F260326-003 LFOAM-1-AC		B7	A0 WATER	1009.82mL 20.00mL	NA	NA	HOAC/ACN	4.5	HOAC/ACN	5.0	S0UL-09GCSV0172
7/01/09	7/13/09	G9F270189-001 LFRD6-1-AC		B7	A0 WATER	985.56mL 20.00mL	NA	NA	HOAC/ACN	4.5	HOAC/ACN	5.0	S0UL-09GCSV0172
7/02/09	7/13/09	G9F270189-002 LFRD7-1-AC		B7	A0 WATER	999.12mL 20.00mL	NA	NA	HOAC/ACN	4.5	HOAC/ACN	5.0	S0UL-09GCSV0172
7/02/09	7/13/09	G9F270189-003 LFRD8-1-AC		B7	A0 WATER	991.73mL 20.00mL	NA	NA	HOAC/ACN	4.5	HOAC/ACN	5.0	S0UL-09GCSV0172
7/02/09	7/13/09	G9F270189-004 LFRD9-1-AC		B7	A0 WATER	996.78mL 20.00mL	NA	NA	HOAC/ACN	4.5	HOAC/ACN	5.0	S0UL-09GCSV0172

RQC058

TestAmerica Laboratories, Inc.  
EXTRACTION BENCH WORKSHEET

Run Date: 7/02/09  
Time: 11:49:39

\*\*\*\*\*  
\* QC BATCH: 9182192 \*  
\* PREP DATE: 7/01/09 9:00 \*  
\* COMP DATE: 7/02/09 10:40 \*  
\*\*\*\*\*

EXTR EXPR	ANL LOT# / DUE WORK ORDER	MSRUN# / TEST FLGS	EXT MTH	MAJRIX	INIT/FIN WT/VOL	PH'S ADJ1	ADJ2	EXTRACTION VOL	SOLVENTS VOL EXCHANGE	VOL	SPIKE STANDARD / SURROGATE ID		
7/02/09	G9F270189-005 LFV9A-1-AC		B7	A0	WATER	985.33mL 20.00mL	NA	NA	HOAC/ACN	4.5	HOAC/ACN	5.0	50UL-09GCSV0172
COMMENTS:													
7/02/09	G9F300242-001 LFV9A-1-AC		B7	A0	WATER	1014.4mL 20.00mL	NA	NA	HOAC/ACN	4.5	HOAC/ACN	5.0	50UL-09GCSV0172
COMMENTS:													
7/02/09	G9F300242-002 LFV9G-1-AC		B7	A0	WATER	975.28mL 20.00mL	NA	NA	HOAC/ACN	4.5	HOAC/ACN	5.0	50UL-09CCSV0172
COMMENTS:													
7/02/09	G9F300242-003 LFV9J-1-AC		B7	A0	WATER	1018.21mL 20.00mL	NA	NA	HOAC/ACN	4.5	HOAC/ACN	5.0	50UL-09GCSV0172
COMMENTS:													
7/03/09	G9F300242-004 LFV9K-1-AC		B7	A0	WATER	1028.11mL 20.00mL	NA	NA	HOAC/ACN	4.5	HOAC/ACN	5.0	50UL-09GCSV0172
COMMENTS:													
7/03/09	G9F300242-004 LFV9K-1-AFS		B7	A0	WATER	1018.28mL 20.00mL	NA	NA	HOAC/ACN	4.5	HOAC/ACN	5.0	SEE BENCH SHEET 50UL-09GCSV0172
COMMENTS:													
7/03/09	G9F300242-004 LFV9K-1-AGD		B7	A0	WATER	1017.3mL 20.00mL	NA	NA	HOAC/ACN	4.5	HOAC/ACN	5.0	SEE BENCH SHEET 50UL-09GCSV0172
COMMENTS:													
7/03/09	G9F300242-005 LFV9M-1-AC		B7	A0	WATER	1016.72mL 20.00mL	NA	NA	HOAC/ACN	4.5	HOAC/ACN	5.0	50UL-09GCSV0172
COMMENTS:													
7/01/09	G9G010000-192 LFW79-1-AAB		B7	A0	WATER	1000mL 20.00mL	NA	NA	HOAC/ACN	4.5	HOAC/ACN	5.0	50UL-09GCSV0172
COMMENTS:													
7/01/09	G9G010000-192 LFW79-1-ACC		B7	A0	WATER	1000mL 20.00mL	NA	NA	HOAC/ACN	4.5	HOAC/ACN	5.0	SEE BENCH SHEET 50UL-09GCSV0172
COMMENTS:													

.1% HOAC.ACN 2991-95F; .1%HOAC/H2O 2991-87B; FILTER MILLIPORE LOT R5PN29184  
SPE COLUMN WATER LOT 00333984A; SODIUM CHLORIDE MALLINKRODT LOT H07582  
LCS, MS/MSD SPIKE 200L-09GCSV0037; 100UL-09GCSV0149; 100UL-09GCSV0039

R = RUSH  
E = EPA 600  
M = CLIENT REQ MS/MSD  
C = CLP  
D = EXP.DEL)

NUMBER OF WORK ORDERS IN BATCH: 17



TestAmerica West Sacramento  
INITIAL CALIBRATION DATA

Start Cal Date : 10-JUN-2009 15:33  
 End Cal Date : 11-JUN-2009 13:47  
 Quant Method : ESTD  
 Origin : Disabled  
 Target Version : 4.14  
 Integrator : Falcon  
 Method file : \\Terastation\share\GCdata\LC10.I\06102009.B\8330AB.M  
 Last Edit : 11-Jun-2009 15:06 kenneyf  
 Curve Type : Average

Calibration File Names:

- Level 1: \\Terastation\share\GCdata\GCdata\LC10.I\06102009.B\A-000002.d
- Level 2: \\Terastation\share\GCdata\GCdata\LC10.I\06102009.B\A-000003.d
- Level 3: \\Terastation\share\GCdata\GCdata\LC10.I\06102009.B\A-000015.d
- Level 4: \\Terastation\share\GCdata\GCdata\LC10.I\06102009.B\A-000005.d
- Level 5: \\Terastation\share\GCdata\GCdata\LC10.I\06102009.B\A-000006.d
- Level 6: \\Terastation\share\GCdata\GCdata\LC10.I\06102009.B\A-000007.d
- Level 7: \\Terastation\share\GCdata\GCdata\LC10.I\06102009.B\A-000008.d
- Level 8: \\Terastation\share\GCdata\GCdata\LC10.I\06102009.B\A-000009.d

Compound	5.000 Level 1	10.000 Level 2	20.000 Level 3	50.000 Level 4	100.000 Level 5	200.000 Level 6	RRF	% RSD
2 HMX	119 ✓ 116	120 ✓ 123	115 ✓	125 ✓	123 ✓	124	121	3.066
3 RDX	81.80000 ✓ 69.16800	75.90000 ✓ 79.81500	77.15000 ✓	80.72000 ✓	78.45000 ✓	80.13000	77.89163	5.156
4 EGDN	++++ ++++	++++ ++++	++++	++++	++++	++++	++++	++++
5 Picric ACID	++++ 74.63400	++++ 74.87800	91.72000 ✓	85.68000 ✓	84.85500 ✓	84.88200	82.77483	8.120
6 1,3,5-Trinitrobenzene	161 136	146 147	145	150	147	148	148	4.610
7 1,3-Dinitrobenzene	145 128	142 141	138	142	139	142	140	3.715
8 3,5-Dinitroaniline	97.20000 82.20800	90.80000 89.25000	89.35000	91.80000	88.83000	90.35000	89.97350	4.581
9 TETRYL	97.60000 ✓ 92.43000	95.10000 ✓ 94.40900	77.40000 ✓	96.02000 ✓	94.31000 ✓	95.06000	92.79113	6.888

*Handwritten:* 6/11/09

TestAmerica West Sacramento

INITIAL CALIBRATION DATA

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 Target Version : 4.14  
 Integrator : Falcon  
 Method file : \\Terastation\share\GCdata\LC10.I\06102009.B\8330AB.M  
 Last Edit : 11-Jun-2009 15:06 kenneyf  
 Curve Type : Average

Compound	5.000 Level 1	10.000 Level 2	20.000 Level 3	50.000 Level 4	100.000 Level 5	200.000 Level 6	RRF	% RSD
10 Nitrobenzene	63.80000 56.02600	62.30000 61.95700	56.55000	61.44000	61.06000	62.26000	60.67412	4.659
11 Nitroglycerin	+++++	+++++	+++++	+++++	+++++	+++++	+++++	+++++
12 2,4,6-Trinitrotoluene	94.80000 84.22200	90.20000 87.17800	81.90000	88.78000	87.04000	87.55500	87.70937	4.401
13 4-AM-2,6-DNT	67.00000 58.65400	63.30000 61.14600	60.85000	63.26000	61.61000	62.67000	62.31125	3.901
14 2-AM-4,6-DNT	74.80000 69.18400	75.80000 72.13500	73.10000	75.54000	73.53000	74.39000	73.55987	2.932
15 2,6-Dinitrotoluene	57.60000 50.43000	53.30000 52.74900	51.90000	53.68000	53.03000	53.21500	53.23800	3.835
16 2,4-Dinitrotoluene	93.00000 81.43200	86.50000 85.05700	82.60000	86.20000	84.72000	85.49000	85.62488	4.031
17 2-Nitrotoluene	40.80000 33.97400	37.40000 35.45000	33.20000	35.64000	35.21000	35.77500	35.93113	6.486
18 4-Nitrotoluene	48.00000 41.67600	46.40000 43.21300	40.75000	43.56000	43.22000	43.56500	43.79800	5.390
19 3-Nitrotoluene	47.00000 41.51600	44.60000 42.82400	39.85000	43.24000	42.93000	43.08000	43.13000	4.852
20 PETN	+++++	+++++	+++++	+++++	+++++	+++++	+++++	+++++



TestAmerica West Sacramento

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 Integrator : Falcon  
 Method file : \\Terastation\share\GCdata\LC10.I\06102009.B\8330AB.M  
 Last Edit : 11-Jun-2009 15:06 kenneyf  
 Curve Type : Average

Compound	5.000 Level 1	10.000 Level 2	20.000 Level 3	50.000 Level 4	100.000 Level 5	200.000 Level 6	RRF	% RSD
	500.000 Level 7	1000.000 Level 8						
\$ 1 3,4-Dinitrotoluene	++++ 46.08000	50.10000 48.38600	46.40000	46.48000	45.87000	46.85500	47.16729	3.251

TestAmerica West Sacramento

INITIAL CALIBRATION DATA

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 Origin : Disabled  
 Target Version : 4.14  
 Integrator : Falcon  
 Method file : \\Terastation\share\GCdata\LC10.I\06102009.B\8330AB.M\83  
 Last Edit : 11-Jun-2009 15:14 kenneyf  
 Curve Type : Average

Calibration File Names:

Level 2: \\Terastation\share\GCdata\GCdata\LC10.I\06102009.B\A-000003.d\A-  
 Level 3: \\Terastation\share\GCdata\GCdata\LC10.I\06102009.B\A-000015.d\A-  
 Level 4: \\Terastation\share\GCdata\GCdata\LC10.I\06102009.B\A-000005.d\A-  
 Level 5: \\Terastation\share\GCdata\GCdata\LC10.I\06102009.B\A-000006.d\A-  
 Level 6: \\Terastation\share\GCdata\GCdata\LC10.I\06102009.B\A-000007.d\A-  
 Level 7: \\Terastation\share\GCdata\GCdata\LC10.I\06102009.B\A-000008.d\A-  
 Level 8: \\Terastation\share\GCdata\GCdata\LC10.I\06102009.B\A-000009.d\A-

Compound	10.000 Level 2	20.000 Level 3	50.000 Level 4	100.000 Level 5	200.000 Level 6	500.000 Level 7	RRF	% RSD
2 HMX	+++++	+++++	+++++	+++++	+++++	+++++	+++++	+++++
3 RDX	+++++	+++++	+++++	+++++	+++++	+++++	+++++	+++++
4 EGDN	+++++	+++++	+++++	+++++	+++++	+++++	+++++	+++++
5 Picric ACID	+++++ 110	134	126	125	125	110	122	8.080
6 1,3,5-Trinitrobenzene	+++++	+++++	+++++	+++++	+++++	+++++	+++++	+++++
7 1,3-Dinitrobenzene	+++++	+++++	+++++	+++++	+++++	+++++	+++++	+++++
8 3,5-Dinitroaniline	+++++	+++++	+++++	+++++	+++++	+++++	+++++	+++++
9 TETRYL	+++++	+++++	+++++	+++++	+++++	+++++	+++++	+++++

TestAmerica West Sacramento

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 Method file : \\Terastation\share\GCdata\LC10.I\06102009.B\8330AB.M\83  
 Last Edit : 11-Jun-2009 15:14 kenneyf  
 Curve Type : Average

Compound	10.000	20.000	50.000	100.000	200.000	500.000	RRF	% RSD
	Level 2	Level 3	Level 4	Level 5	Level 6	Level 7		
	1000.000							
	Level 8							
10 Nitrobenzene	+++++	+++++	+++++	+++++	+++++	+++++	+++++	+++++
11 Nitroglycerin	+++++	61.28000	59.82000	60.58000	61.86500	59.75400	60.83467	1.527
12 2,4,6-Trinitrotoluene	+++++	+++++	+++++	+++++	+++++	+++++	+++++	+++++
13 4-AM-2,6-DNT	+++++	+++++	+++++	+++++	+++++	+++++	+++++	+++++
14 2-AM-4,6-DNT	+++++	+++++	+++++	+++++	+++++	+++++	+++++	+++++
15 2,6-Dinitrotoluene	+++++	+++++	+++++	+++++	+++++	+++++	+++++	+++++
16 2,4-Dinitrotoluene	+++++	+++++	+++++	+++++	+++++	+++++	+++++	+++++
17 2-Nitrotoluene	+++++	+++++	+++++	+++++	+++++	+++++	+++++	+++++
18 4-Nitrotoluene	+++++	+++++	+++++	+++++	+++++	+++++	+++++	+++++
19 3-Nitrotoluene	+++++	+++++	+++++	+++++	+++++	+++++	+++++	+++++
20 PETN	+++++	32.65000	30.36000	31.91000	32.25000	32.38000	32.05033	2.748

TestAmerica West Sacramento

INITIAL CALIBRATION DATA

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 Target Version : 4.14  
 Integrator : Falcon  
 Method file : \\Terastation\share\GCdata\LC10.I\06102009.B\8330AB.M\83  
 Last Edit : 11-Jun-2009 15:14 kenneyf  
 Curve Type : Average

Compound	10.000 Level 2	20.000 Level 3	50.000 Level 4	100.000 Level 5	200.000 Level 6	500.000 Level 7	RRF	% RSD
	1000.000 Level 8							
\$ 1,3,4-Dinitrotoluene	* +++++ 93.14200	90.40000	88.76000	88.28000	90.51000	88.03667	89.85478	2.142

*\* Level 2 dropped due to poor integration  
 JMK  
 6/12/09*

Chromatography Summary

Injection Date: 6/10/2009 23:09 Operator: fhk  
 DataFile: IC10.IV05102009.DVA-000011.D Vial Num: 20  
 Instrument ID: IC10

Method 8330 Target Analyte Results

Sample: **ICV\_6 08GCSV0397**  
**100/200/100/100ng/mL**

Method File: IC10.IV05102009.BW8330AB.M  
 Start Cal Date: 6/10/2009 15:33 End Cal Date: 6/11/2009 13:47

Matrix: NONE SubList: CAL.sub SpikeList:  
 Samp. Info: ICV\_6 08GCSV0397 100/200/100/100ng/mL?  
 Misc. Info: ;6;;;3;CAL;sub;:0;1

Dilution Factor	Extract Volume	Sample Volume	Sample Weight
1X	0 mL	0 mL	0 g

Compound Name	RT	Signal 1 UV 250-265					Flag	Signal 2 UV 358-205					Limits(%)	Flag
		Response	PPB	Spike Level	%D	Result		RT	Response	PPB	Spike Level	%D		
<b>3,4-Dinitrotoluene</b>				200	-100%	Fails				200	-100%	Fails	(±15)	
HMX	5.29	26153	216.4000<	200	8%	Acceptable				200	-100%	Fails	(±15)	45
RDX	7.84	16084	206.5000<	200	3%	Acceptable				200	-100%	Fails	(±15)	45
Picric ACID	9.00	41754	504.4000 ✓	500	1%	Acceptable	9.00	61509	505.5000< ✓	500	1%	Acceptable	(±15)	
1,3,5-Trinitrobenzene	10.20	29097	197.1000<	200	-1%	Acceptable				200	-100%	Fails	(±15)	45
1,3-Dinitrobenzene	13.25	27834	199.2000<	200	0%	Acceptable				200	-100%	Fails	(±15)	45
TETRYL	14.30	17776	191.6000<	200	-4%	Acceptable				200	-100%	Fails	(±15)	45
Nitrobenzene	15.16	12143	200.1000<	200	0%	Acceptable				200	-100%	Fails	(±15)	45
2,4,6-Trinitrotoluene	16.53	17352	197.8000<	200	-1%	Acceptable				200	-100%	Fails	(±15)	45
4-AM-2,6-DNT	17.31	11969	192.1000< ✓	200	-4%	Acceptable				200	-100%	Fails	(±15)	45
2-AM-4,6-DNT	18.57	13851	188.3000<	200	-6%	Acceptable				200	-100%	Fails	(±15)	45
2,6-Dinitrotoluene	20.28	10445	196.2000<	200	-2%	Acceptable				200	-100%	Fails	(±15)	45
2,4-Dinitrotoluene	21.00	16653	194.5000<	200	-3%	Acceptable				200	-100%	Fails	(±15)	45
2-Nitrotoluene	24.68	7069	196.7000<	200	-2%	Acceptable				200	-100%	Fails	(±15)	45
4-Nitrotoluene	26.50	8526	194.7000<	200	-3%	Acceptable				200	-100%	Fails	(±15)	45
3-Nitrotoluene	28.49	8451	195.9000< ✓	200	-2%	Acceptable				200	-100%	Fails	(±15)	45
<b>Nitroglycerin</b>				200	-100%	Fails	15.78	12548	206.3000< ✓	200	3%	Acceptable	(±15)	45
PETN				200	-100%	Fails	30.10	5830	181.9000< ✓	200	-9%	Acceptable	(±15)	45
3,5-Dinitroaniline	13.99	18460	205.2000<	200	3%	Acceptable				200	-100%	Fails	(±15)	45
<b>EGDN</b>				200	-100%	Fails				200	-100%	Fails	(±15)	

Notes: M = Manually Integrated Signals Differ by More Than 40%  
 D = Operator Disabled Result Signals Differ by More Than 50%  
 O = Over Calibration Range

TestAmerica West Sacramento  
GC/LC INSTRUMENT LOG

Inst ID: LC10                                      Batch ID: 07022009  
Method : Method 8330                              Test : SOP SAC-LC-0009  
ICAL Date: See Calibration Report

Date	Time	Operator	Sample ID	File ID	Vol or wt	Extract Vol	Diln	Comments
02-JUL-2009	15:27	fhk	Primer	A-000001.	0 g	0 mL	1	
02-JUL-2009	16:18	fhk	Primer	A-000002.	0 g	0 mL	1	
02-JUL-2009	17:09	fhk	STD_6 09GCSV0054 .4K/.2/.2/.2/	A-000003.	0 g	0 mL	1	
02-JUL-2009	18:00	fhk	LFW791AA 9182192 G9G010000-192	A-000004.	1000 mL	20 mL	1	
02-JUL-2009	18:50	fhk	LFW791AC 9182192 G9G010000-192	A-000005.	1000 mL	20 mL	1	
02-JUL-2009	19:40	fhk	LFQAKIAC 9182192 G9F260326-1 1	A-000006.	997.78 mL	20 mL	1	
02-JUL-2009	20:31	fhk	LFQALIAC 9182192 G9F260326-2 1	A-000007.	991.12 mL	20 mL	1	
02-JUL-2009	21:22	fhk	LFQAMIAC 9182192 G9F260326-3 1	A-000008.	1009.82 mL	20 mL	1	
02-JUL-2009	22:13	fhk	LFRD6IAC 9182192 G9F270189-1 1	A-000009.	985.56 mL	20 mL	1	
02-JUL-2009	23:03	fhk	LFRD7IAC 9182192 G9F270189-2 1	A-000010.	999.12 mL	20 mL	1	
02-JUL-2009	23:54	fhk	LFRD8IAC 9182192 G9F270189-3 1	A-000011.	991.73 mL	20 mL	1	
03-JUL-2009	00:45	fhk	STD_5 09GCSV0053 2K/.1/.1/.1/	A-000012.	0 g	0 mL	1	
03-JUL-2009	01:36	fhk	LFRD9IAC 9182192 G9F270189-4 1	A-000013.	996.78 mL	20 mL	1	
03-JUL-2009	02:27	fhk	LFREA1AC 9182192 G9F270189-5 1	A-000014.	985.33 mL	20 mL	1	
03-JUL-2009	03:17	fhk	LFV9A1AC 9182192 G9F300242-1 1	A-000015.	1014.4 mL	20 mL	1	
03-JUL-2009	04:07	fhk	LFV9G1AC 9182192 G9F300242-2 1	A-000016.	975.28 mL	20 mL	1	
03-JUL-2009	04:58	fhk	LFV9J1AC 9182192 G9F300242-3 1	A-000017.	1018.21 mL	20 mL	1	
03-JUL-2009	05:49	fhk	LFV9K1AC 9182192 G9F300242-4 1	A-000018.	1028.11 mL	20 mL	1	
03-JUL-2009	06:39	fhk	MS LFV9K1AF 9182192 G9F300242-4 S	A-000019.	1018.28 mL	20 mL	1	
03-JUL-2009	07:30	fhk	MSD LFV9K1AG 9182192 G9F300242-4 D	A-000020.	1017.3 mL	20 mL	1	
03-JUL-2009	08:20	fhk	LFV9M1AC 9182192 G9F300242-5 1	A-000021.	1016.72 mL	20 mL	1	
03-JUL-2009	09:11	fhk	STD_6 09GCSV0054 .4K/.2/.2/.2/	A-000022.	0 g	0 mL	1	
03-JUL-2009	10:01	fhk	09GCSV0234 LEVEL 1	A-000023.	0 g	0 mL	1	
03-JUL-2009	10:52	fhk	09GCSV0235 LEVEL 2	A-000024.	0 g	0 mL	1	
03-JUL-2009	11:42	fhk	09GCSV0236 LEVEL 3	A-000025.	0 g	0 mL	1	
03-JUL-2009	12:33	fhk	09GCSV0237 LEVEL 4	A-000026.	0 g	0 mL	1	
03-JUL-2009	13:23	fhk	09GCSV0238 LEVEL 5	A-000027.	0 g	0 mL	1	
03-JUL-2009	14:14	fhk	09GCSV0239 LEVEL 5	A-000028.	0 g	0 mL	1	
03-JUL-2009	15:04	fhk	09GCSV0240 LEVEL 6	A-000029.	0 g	0 mL	1	
03-JUL-2009	15:55	fhk	09GCSV0241 LEVEL 7	A-000030.	0 g	0 mL	1	
03-JUL-2009	16:45	fhk	09GCSV0242 LEVEL 8	A-000031.	0 g	0 mL	1	
03-JUL-2009	17:36	fhk	09GCSV0243 MRL	A-000032.	0 g	0 mL	1	

Chromatography Summary

Injection Date: 7/2/2009 17:09 Operator: fhk  
 DataFile: LC10.N07022009.BVA-000003.D Vial Num: 2  
 Instrument ID: LC10

Method 8330 Target Analyte Results

Sample: **STD\_6 09GCSV0054 4K/2/2/2/2**

Method File: LC10.N07022009.D\8330AB.M  
 Start Cal Date: 6/10/2009 15:33 End Cal Date: 6/11/2009 13:47

Matrix: NONE SubList: CAL sub SpikeList:  
 Samp. Info: STD\_6 09GCSV0054 4K/2/2/2/2  
 Misc. Info: ,6, , , ;3;CAL.sub; ;0;1

Dilution Factor	Extract Volume	Sample Volume	Sample Weight
1X	0 mL	0 mL	0 g

Compound Name	RT	Signal 1 UV 250-265					Signal 2 UV 358-205					Limits(%)	Flag			
		Response	PPB	Spike Level	%D	Result	Flag	RT	Response	PPB	Spike Level			%D	Result	Flag
3,4-Dinitrotoluene	18.20	9636	204	3000<	200	2%	Acceptable	18.20	18445	213	4000	200	7%	Acceptable	(±15)	
HMX	5.29	25432	210	5000<	200	5%	Acceptable					200	-100%	Fails	(±15)	45
RDX	7.85	16295	209	2000<	200	5%	Acceptable					200	-100%	Fails	(±15)	45
Picric ACID	8.99	41811	505	1000	500	1%	Acceptable	8.99	61520	505	6000<	500	1%	Acceptable	(±15)	
1,3,5-Trinitrobenzene	10.20	30367	205	7000<	200	3%	Acceptable					200	-100%	Fails	(±15)	45
1,3-Dinitrobenzene	13.26	29099	208	3000<	200	4%	Acceptable					200	-100%	Fails	(±15)	45
TETRYL	14.32	19114	206	0000<	200	3%	Acceptable					200	-100%	Fails	(±15)	45
Nitrobenzene	15.19	12872	212	1000<	200	6%	Acceptable					200	-100%	Fails	(±15)	45
2,4,6-Trinitrotoluene	16.55	18075	206	1000<	200	3%	Acceptable					200	-100%	Fails	(±15)	45
4-AM-2,6-DNT	17.55	12873	206	6000<	200	3%	Acceptable					200	-100%	Fails	(±15)	45
2-AM-4,6-DNT	18.61	15046	204	5000<	200	2%	Acceptable					200	-100%	Fails	(±15)	45
2,6-Dinitrotoluene	20.34	10970	206	0000<	200	3%	Acceptable					200	-100%	Fails	(±15)	45
2,4-Dinitrotoluene	21.06	17567	205	2000<	200	3%	Acceptable					200	-100%	Fails	(±15)	45
2-Nitrotoluene	24.79	7368	205	0000<	200	3%	Acceptable					200	-100%	Fails	(±15)	45
4-Nitrotoluene	26.61	3982	205	1000<	200	3%	Acceptable					200	-100%	Fails	(±15)	45
3-Nitrotoluene	28.64	3962	207	8000<	200	4%	Acceptable					200	-100%	Fails	(±15)	45
Nitroglycerin					200	-100%	Fails	15.81	12531	206	0000<	200	3%	Acceptable	(±15)	45
PETN					200	-100%	Fails	30.22	6862	214	1000<	200	7%	Acceptable	(±15)	45
3,5-Dinitroaniline	14.01	18502	205	6000<	200	3%	Acceptable					200	-100%	Fails	(±15)	45
EGDN					200	-100%	Fails					200	-100%	Fails	(±15)	

Notes: M = Manually Integrated Signals Differ by More Than 40%  
 D = Operator Disabled Result Signals Differ by More Than 50%  
 O = Over Calibration Range

Chromatography Summary

Injection Date: 7/3/2009 0:45 Operator: fhk  
 Data File: LC10.I07022009.BVA-000012.D Vial Num: 3  
 Instrument ID: LC10

Method 8330 Target Analyte Results

Sample: **STD\_5 09GCSV0053 .2K/1/1/1/1**

Method File: LC10 I07022009.BV8330AB.M  
 Start Cal Date: 6/10/2009 15:33 End Cal Date: 6/11/2009 13:47

Matrix: NONE SubList: CAL sub SpikeList:  
 Samp. Info: STD\_5 09GCSV0053 2K/1/1/1/1,2  
 Misc. Info: ;S; ; ;3;CAL.sub; ;0;1

Dilution Factor	Extract Volume	Sample Volume	Sample Weight
1X	0 mL	0 mL	0 g

Signal 1 UV 250-265								Signal 2 UV 358-205								
Compound Name	RT	Response	PPB	Spike Level	%D	Result	Flag	RT	Response	PPB	Spike Level	%D	Result	Flag	Limits(%)	Flag
3,4-Dinitrotoluene	18.21	4770	101 1000<	100	1%	Acceptable		18.21	9105	105 3000	100	5%	Acceptable		(±15)	
HMX	5.29	12564	104 8000<	100	5%	Acceptable					100	-100%	Fails		(±15)	45
RDX	7.86	8028	103 1000<	100	3%	Acceptable					100	-100%	Fails		(±15)	45
Picric ACID	9.05	16980	205 1000	200	3%	Acceptable		9.05	25011	205 5000<	200	3%	Acceptable		(±15)	
1,3,5-Trinitrobenzene	10.21	15116	102 4000<	100	2%	Acceptable					100	-100%	Fails		(±15)	45
1,3-Dinitrobenzene	13.27	14453	101 5000<	100	4%	Acceptable					100	-100%	Fails		(±15)	45
TETRYL	14.34	9510	102 5000<	100	3%	Acceptable					100	-100%	Fails		(±15)	45
Nitrobenzene	15.20	6322	104 2000<	100	4%	Acceptable					100	-100%	Fails		(±15)	45
2,4,6-Trinitrotoluene	16.57	9018	102 8000<	100	3%	Acceptable					100	-100%	Fails		(±15)	45
4-AM-2,6-DNT	17.56	6403	102 8000<	100	3%	Acceptable					100	-100%	Fails		(±15)	45
2-AM-4,6-DNT	18.63	7508	102 1000<	100	2%	Acceptable					100	-100%	Fails		(±15)	45
2,6-Dinitrotoluene	20.35	5455	102 5000<	100	3%	Acceptable					100	-100%	Fails		(±15)	45
2,4-Dinitrotoluene	21.08	8726	101 9000<	100	2%	Acceptable					100	-100%	Fails		(±15)	45
2-Nitrotoluene	24.79	3635	101 2000<	100	1%	Acceptable					100	-100%	Fails		(±15)	45
4-Nitrotoluene	26.61	4448	101 6000<	100	2%	Acceptable					100	-100%	Fails		(±15)	45
3-Nitrotoluene	28.63	4394	101 9000<	100	2%	Acceptable					100	-100%	Fails		(±15)	45
Nitroglycerin				100	-100%	Fails		15.82	6254	102 8000<	100	3%	Acceptable		(±15)	45
PETN				100	-100%	Fails		30.25	3304	103 1000<	100	3%	Acceptable		(±15)	45
3,5-Dinitroaniline	14.02	9179	102 0000<	100	2%	Acceptable					100	-100%	Fails		(±15)	45
EGDN				100	-100%	Fails					100	-100%	Fails		(±15)	

Notes: M = Manually Integrated Signals Differ by More Than 40%  
 D = Operator Disabled Result Signals Differ by More Than 50%  
 O = Over Calibration Range



Chromatography Summary

Injection Date: 7/3/2009 9:11 Operator: fhk  
 DataFile: LC10 R07022009.BA-000022.D Vial Num: 4  
 Instrument ID: LC10

Method 8330 Target Analyte Results

Sample: **STD\_6 09GCSV0054 .4K/2/2/2/2**

Method File: LC10 R07022009.BA8330AB.M  
 Start Cal Date: 6/10/2009 15:33 End Cal Date: 6/11/2009 13:47

Matrix: NONE SubList: CAL.sub SpikeList:  
 Samp. Info: STD\_6 09GCSV0054 .4K/2/2/2/2  
 Misc. Info: .6, ; ; 3,CAL.sub, .0,1

Dilution Factor	Extract Volume	Sample Volume	Sample Weight
1X	0 mL	0 mL	0 g

Compound Name	Signal 1 UV 250-265							Signal 2 UV 358-205								
	RT	Response	PPB	Spike Level	%D	Result	Flag	RT	Response	PPB	Spike Level	%D	Result	Flag	Limits(%)	Flag
3,4-Dinitrotoluene	18.24	9589	203.3000<	200	2%	Acceptable		18.23	18290	211.6000	200	6%	Acceptable		(±15)	
HMX	5.30	25440	210.6000<	200	5%	Acceptable					200	-100%	Fails		(±15)	45
RDX	7.87	16268	208.8000<	200	4%	Acceptable					200	-100%	Fails		(±15)	45
Picric ACID	9.04	42655	515.3000	500	3%	Acceptable		9.04	62723	515.5000<	500	3%	Acceptable		(±15)	
1,3,5-Trinitrobenzene	10.22	30282	205.1000<	200	3%	Acceptable					200	-100%	Fails		(±15)	45
1,3-Dinitrobenzene	13.29	28991	207.5000<	200	4%	Acceptable					200	-100%	Fails		(±15)	45
TETRYL	14.35	19191	206.8000<	200	3%	Acceptable					200	-100%	Fails		(±15)	45
Nitrobenzene	15.21	12738	209.9000<	200	5%	Acceptable					200	-100%	Fails		(±15)	45
2,4,6-Trinitrotoluene	16.58	18022	205.5000<	200	3%	Acceptable					200	-100%	Fails		(±15)	45
4-AM-2,6-DNT	17.58	12828	205.9000<	200	3%	Acceptable					200	-100%	Fails		(±15)	45
2-AM-4,6-DNT	18.65	15059	204.7000<	200	2%	Acceptable					200	-100%	Fails		(±15)	45
2,6-Dinitrotoluene	20.38	10882	204.4000<	200	2%	Acceptable					200	-100%	Fails		(±15)	45
2,4-Dinitrotoluene	21.10	17483	204.2000<	200	2%	Acceptable					200	-100%	Fails		(±15)	45
2-Nitrotoluene	24.82	7294	203.0000<	200	2%	Acceptable					200	-100%	Fails		(±15)	45
4-Nitrotoluene	26.64	8876	202.6000<	200	1%	Acceptable					200	-100%	Fails		(±15)	45
3-Nitrotoluene	28.67	8836	204.9000<	200	2%	Acceptable					200	-100%	Fails		(±15)	45
Nitroglycerin				200	-100%	Fails		15.84	12481	205.2000<	200	3%	Acceptable		(±15)	45
PETN				200	-100%	Fails		30.27	6681	208.4000<	200	4%	Acceptable		(±15)	45
3,5-Dinitroaniline	14.03	18458	205.1000<	200	3%	Acceptable					200	-100%	Fails		(±15)	45
EGDN				200	-100%	Fails					200	-100%	Fails		(±15)	

Notes: M = Manually Integrated Signals Differ by More Than 40%  
 D = Operator Disabled Result Signals Differ by More Than 50%  
 O = Over Calibration Range



TestAmerica West Sacramento

INITIAL CALIBRATION DATA

Start Cal Date : 09-JUN-2009 20:53  
 End Cal Date : 10-JUN-2009 04:20  
 Quant Method : ESTD  
 Origin : Disabled  
 Target Version : 4.14  
 Integrator : HP Genie  
 Method file : \\Terastation\share\GCdata\LC9.I\06092009.B\8330CNAB.M  
 Last Edit : 10-Jun-2009 10:24 kenneyf  
 Curve Type : Average

Calibration File Names:

- Level 1: \\Terastation\share\GCdata\GCdata\LC9.I\06092009.B\C-000004.d
- Level 2: \\Terastation\share\GCdata\GCdata\LC9.I\06092009.B\C-000005.d
- Level 3: \\Terastation\share\GCdata\GCdata\LC9.I\06092009.B\C-000006.d
- Level 4: \\Terastation\share\GCdata\GCdata\LC9.I\06092009.B\C-000007.d
- Level 5: \\Terastation\share\GCdata\GCdata\LC9.I\06092009.B\C-000008.d
- Level 6: \\Terastation\share\GCdata\GCdata\LC9.I\06092009.B\C-000009.d
- Level 7: \\Terastation\share\GCdata\GCdata\LC9.I\06092009.B\C-000010.d
- Level 8: \\Terastation\share\GCdata\GCdata\LC9.I\06092009.B\C-000011.d

Compound	5.000	10.000	20.000	50.000	100.000	200.000	RRF	% RSD
	Level 1	Level 2	Level 3	Level 4	Level 5	Level 6		
2 HMX	49.60000 37.41800	43.70000 39.22000	40.45000	40.46000	39.51000	39.31500	41.20913	9.288
3 RDX	45.40000 38.11800	43.70000 42.45200	42.75000	43.30000	42.16000	42.58500	42.55812	4.852
4 EGDN	+++++	+++++	+++++	+++++	+++++	+++++	+++++	+++++
5 Picric ACID	+++++	+++++	+++++	+++++	+++++	+++++	+++++	+++++
6 1,3,5-Trinitrobenzene	66.80000 63.46600	67.20000 67.27500	67.60000	68.82000	67.02000	67.61500	66.97450	2.306
7 1,3-Dinitrobenzene	95.20000 79.37000	88.70000 88.38800	87.25000	89.40000	87.58000	88.56500	88.05662	4.893
8 3,5-Dinitroaniline	71.20000 60.97200	68.50000 65.89900	68.65000	68.54000	66.12000	66.97500	67.10700	4.478
9 TETRYL	93.20000 85.04400	86.50000 86.33600	85.70000	86.92000	86.05000	86.26500	87.00187	2.950

TestAmerica West Sacramento

INITIAL CALIBRATION DATA

Start Cal Date : 09-JUN-2009 20:53  
 End Cal Date : 10-JUN-2009 04:20  
 Quant Method : ESTD  
 Origin : Disabled  
 Target Version : 4.14  
 Integrator : HP Genie  
 Method file : \\Terastation\share\GCdata\LC9.I\06092009.B\8330CNAB.M  
 Last Edit : 10-Jun-2009 10:24 kenneyf  
 Curve Type : Average

Compound	5.000 Level 1	10.000 Level 2	20.000 Level 3	50.000 Level 4	100.000 Level 5	200.000 Level 6	RRP	% RSD
10 Nitrobenzene	39.40000 35.35200	42.00000 40.30500	38.65000	40.64000	39.79000	40.66500	39.60025	5.005
11 Nitroglycerin	++++ ++++	++++ ++++	++++	++++	++++	++++	++++	++++
12 2,4,6-Trinitrotoluene	104 71.97000	91.00000 73.60800	81.00000	77.16000	74.32000	73.90500	80.89538	13.892
13 4-AM-2,6-DNT	68.80000 64.80800	69.80000 65.14700	68.05000	67.74000	65.37000	65.85500	66.94625	2.818
14 2-AM-4,6-DNT	80.20000 69.79400	76.80000 72.19800	75.45000	74.54000	72.25000	73.07500	74.28838	4.343
15 2,6-Dinitrotoluene	54.80000 45.69200	52.40000 49.41100	49.70000	49.88000	49.11000	49.44000	50.05412	5.280
16 2,4-Dinitrotoluene	83.00000 73.10400	82.40000 78.73900	79.60000	80.22000	78.86000	79.32000	79.40538	3.779
17 2-Nitrotoluene	24.50000 22.18600	25.30000 24.28200	23.52500	24.39000	24.32500	24.51250	24.12756	3.813
18 4-Nitrotoluene	++++ ++++	++++ ++++	++++	++++	++++	++++	++++	++++
19 3-Nitrotoluene	35.00000 28.82000	32.50000 30.96400	30.55000	31.42000	30.76000	31.17500	31.39862	5.672
20 PETN	++++ ++++	++++ ++++	++++	++++	++++	++++	++++	++++

TestAmerica West Sacramento

INITIAL CALIBRATION DATA

Start Cal Date : 09-JUN-2009 20:53  
 End Cal Date : 10-JUN-2009 04:20  
 Quant Method : ESTD  
 Origin : Disabled  
 Target Version : 4.14  
 Integrator : HP Genie  
 Method file : \\Terastation\share\GCdata\LC9.I\06092009.B\8330CNAB.M  
 Last Edit : 10-Jun-2009 10:24 kenneyf  
 Curve Type : Average

Compound	5.000 Level 1	10.000 Level 2	20.000 Level 3	50.000 Level 4	100.000 Level 5	200.000 Level 6	RRF	% RSD
	500.000 Level 7	1000.000 Level 8						
\$ 1 3,4-Dinitrotoluene	+++++	46.40000	40.85000	39.44000	38.04000	38.34500	39.65376	8.405
	35.87333	38.62800						

TestAmerica West Sacramento

INITIAL CALIBRATION DATA

Start Cal Date : 09-JUN-2009 21:56  
 End Cal Date : 10-JUN-2009 04:20  
 Quant Method : ESTD  
 Origin : Disabled  
 Target Version : 4.14  
 Integrator : HP Genie  
 Method file : \\Terastation\share\GCdata\LC9.I\06092009.B\8330CNAB.M\8  
 Last Edit : 10-Jun-2009 10:25 kenneyf  
 Curve Type : Average

Calibration File Names:

Level 2: \\Terastation\share\GCdata\GCdata\LC9.I\06092009.B\C-000005.d\C-0  
 Level 3: \\Terastation\share\GCdata\GCdata\LC9.I\06092009.B\C-000006.d\C-0  
 Level 4: \\Terastation\share\GCdata\GCdata\LC9.I\06092009.B\C-000007.d\C-0  
 Level 5: \\Terastation\share\GCdata\GCdata\LC9.I\06092009.B\C-000008.d\C-0  
 Level 6: \\Terastation\share\GCdata\GCdata\LC9.I\06092009.B\C-000009.d\C-0  
 Level 7: \\Terastation\share\GCdata\GCdata\LC9.I\06092009.B\C-000010.d\C-0  
 Level 8: \\Terastation\share\GCdata\GCdata\LC9.I\06092009.B\C-000011.d\C-0

Compound	10.000 Level 2	20.000 Level 3	50.000 Level 4	100.000 Level 5	200.000 Level 6	500.000 Level 7	RRF	% RSD
2 HMX	+++++	+++++	+++++	+++++	+++++	+++++	+++++	+++++
3 RDX	+++++	+++++	+++++	+++++	+++++	+++++	+++++	+++++
4 EGDN	+++++ 42.86600	39.00000	42.36000	44.25000	44.94000	40.36600	42.29700	5.363
5 Picric ACID	+++++	+++++	+++++	+++++	+++++	+++++	+++++	+++++  <-
6 1,3,5-Trinitrobenzene	+++++	+++++	+++++	+++++	+++++	+++++	+++++	+++++
7 1,3-Dinitrobenzene	+++++	+++++	+++++	+++++	+++++	+++++	+++++	+++++
8 3,5-Dinitroaniline	+++++	+++++	+++++	+++++	+++++	+++++	+++++	+++++
9 TETRYL	+++++	+++++	+++++	+++++	+++++	+++++	+++++	+++++

TestAmerica West Sacramento

INITIAL CALIBRATION DATA

Start Cal Date : 09-JUN-2009 21:56  
 End Cal Date : 10-JUN-2009 04:20  
 Quant Method : ESTD  
 Origin : Disabled  
 Target Version : 4.14  
 Integrator : HP Genie  
 Method file : \\Terastation\share\GCdata\LC9.I\06092009.B\8330CNAB.M\8  
 Last Edit : 10-Jun-2009 10:25 kenneyf  
 Curve Type : Average

Compound	10.000 Level 2	20.000 Level 3	50.000 Level 4	100.000 Level 5	200.000 Level 6	500.000 Level 7	1000.000 Level 8	RRF	% RSD
10 Nitrobenzene	++++	++++	++++	++++	++++	++++		++++	++++
11 Nitroglycerin	++++ 57.30100	55.25000	55.16000	56.31000	57.02000	56.37800		56.23650	1.571
12 2,4,6-Trinitrotoluene	++++	++++	++++	++++	++++	++++		++++	++++
13 4-AM-2,6-DNT	++++	++++	++++	++++	++++	++++		++++	++++
14 2-AM-4,6-DNT	++++	++++	++++	++++	++++	++++		++++	++++
15 2,6-Dinitrotoluene	++++	++++	++++	++++	++++	++++		++++	++++
16 2,4-Dinitrotoluene	++++	++++	++++	++++	++++	++++		++++	++++
17 2-Nitrotoluene	++++	++++	++++	++++	++++	++++		++++	++++
18 4-Nitrotoluene	++++	++++	++++	++++	++++	++++		++++	++++
19 3-Nitrotoluene	++++	++++	++++	++++	++++	++++		++++	++++
20 PETN	++++ 105	104	101	102	104	104		103	1.208

TestAmerica West Sacramento

INITIAL CALIBRATION DATA

Start Cal Date : 09-JUN-2009 21:56  
 End Cal Date : 10-JUN-2009 04:20  
 Quant Method : ESTD  
 Origin : Disabled  
 Target Version : 4.14  
 Integrator : HP Genie  
 Method file : \\Terastation\share\GCdata\LC9.I\06092009.B\8330CNAB.M\8  
 Last Edit : 10-Jun-2009 10:25 kenneyf  
 Curve Type : Average

Compound	10.000 Level 2	20.000 Level 3	50.000 Level 4	100.000 Level 5	200.000 Level 6	500.000 Level 7	RRF	% RSD
	1000.000 Level 8							
\$ 1,3,4-Dinitrotoluene	+++++	+++++	+++++	+++++	+++++	+++++	+++++	+++++



Chromatography Summary

Injection Date: 6/10/2009 6:29 Operator: fhk  
 DataFile: LC9 I06092009 BVC-000013.D Vial Num: 20  
 Instrument ID: LC9

Method 8330 Target Analyte Results

Method File: LC9 I06092009 B8330CNAB.M  
 Start Cal Date: 6/9/2009 20:53 End Cal Date: 6/10/2009 4:20

Sample: **ICV\_6 08GCSV0397**  
**200/500/200/200/200**

Dilution Factor	Extract Volume	Sample Volume	Sample Weight
1X	0 mL	0 mL	0 g

Matrix: NONE SubList: CAL.sub SpikeList:  
 Samp. Info: ICV\_6 08GCSV0397 200/500/200/200/200;2  
 Misc. Info: ; 6, ., ; 3; CAL.sub; ; 0; 1

Signal 1 UV 250-265								Signal 2 UV 358-205								
Compound Name	RT	Response	PPB	Spike Level	%D	Result	Flag	RT	Response	PPB	Spike Level	%D	Result	Flag	Limits(%)	Flag
3,4-Dinitrotoluene	31.18	259	6.5320<	200	-97%	Fails	Not in list				200	-100%	Fails		(±15)	45
HMX	35.11	8323	202.0000<	200	1%	Acceptable					200	-100%	Fails		(±15)	45
RDX	25.60	8556	201.0000<	200	1%	Acceptable					200	-100%	Fails		(±15)	45
Picric ACID				500	-100%	Fails					500	-100%	Fails		(±15)	
1,3,5-Trinitrobenzene	21.78	13318	198.8000<	200	-1%	Acceptable					200	-100%	Fails		(±15)	45
1,3-Dinitrobenzene	19.50	17403	197.6000<	200	-1%	Acceptable					200	-100%	Fails		(±15)	45
TETRYL	40.80	16134	185.4000<	200	-7%	Acceptable					200	-100%	Fails		(±15)	45
Nitrobenzene	16.81	7967	201.2000<	200	1%	Acceptable					200	-100%	Fails		(±15)	45
2,4,6-Trinitrotoluene	32.92	14562	180.0000<	200	-10%	Acceptable					200	-100%	Fails		(±15)	45
4-AM-2,6-DNT	29.55	12516	187.0000<	200	-7%	Acceptable					200	-100%	Fails		(±15)	45
2-AM-4,6-DNT	29.21	14216	191.4000<	200	-4%	Acceptable					200	-100%	Fails		(±15)	45
2,6-Dinitrotoluene	27.26	9755	194.9000<	200	-3%	Acceptable					200	-100%	Fails		(±15)	45
2,4-Dinitrotoluene	26.42	15486	195.0000<	200	-3%	Acceptable					200	-100%	Fails		(±15)	45
2-Nitrotoluene	22.83	9684	401.4000<	400	0%	Acceptable					400	-100%	Fails		(±15)	45
4-Nitrotoluene				400	-100%	Fails					400	-100%	Fails		(±15)	
3-Nitrotoluene	23.35	6108	194.5000<	200	-3%	Acceptable					200	-100%	Fails		(±15)	45
Nitroglycerin				200	-100%	Fails		38.67	11785	209.6000<	200	5%	Acceptable		(±15)	45
PETN				200	-100%	Fails		49.13	19084	184.7000<	200	-8%	Acceptable		(±15)	45
3,5-Dinitroaniline	25.06	13705	204.2000<	200	2%	Acceptable					200	-100%	Fails		(±15)	45
EGDN				200	-100%	Fails		18.01	9041	213.8000<	200	7%	Acceptable		(±15)	45

Notes: M = Manually Integrated Signals Differ by More Than 40%  
 D = Operator Disabled Result Signals Differ by More Than 50%  
 O = Over Calibration Range



Chromatography Summary

Injection Date: 7/11/2009 5:54 Operator: fhk  
 DataFile: LC9\_I07102009A\_BAC-000013.D Vial Num: 3  
 Instrument ID: LC9

Method 8330 Target Analyte Results

Sample: **STD\_6 09GCSV0054 .5K/2/2/2/2**

Method File: LC9\_I07102009A\_BA8330CNAB.M  
 Start Cal Date: 6/9/2009 20:53 End Cal Date: 6/10/2009 4:20

Matrix: NONE SubList: CAL.sub SpikeList:  
 Samp. Info: STD\_6 09GCSV0054 .5K/2/2/2/2,2  
 Misc. Info: ; 6; ; ; 3; CAL.sub, ; 0; 1; 0

Dilution Factor	Extract Volume	Sample Volume	Sample Weight
1X	0 mL	0 mL	0 g

Signal 1 UV 250-265								Signal 2 UV 358-205								
Compound Name	RT	Response	PPB	Spike Level	%D	Result	Flag	RT	Response	PPB	Spike Level	%D	Result	Flag	Limits(%)	Flag
3,4-Dinitrotoluene	30.69	8487	214.0000<	200	7%	Acceptable					200	-100%	Fails		(±15)	45
HMX	34.67	7594	184.3000<	200	-8%	Acceptable					200	-100%	Fails		(±15)	45
RDX	25.16	8673	203.8000<	200	2%	Acceptable					200	-100%	Fails		(±15)	45
<b>Picric ACID</b>				<b>500</b>	<b>-100%</b>	<b>Fails</b>					<b>500</b>	<b>-100%</b>	<b>Fails</b>		<b>(±15)</b>	
1,3,5-Trinitrobenzene	21.54	13992	208.9000<	200	4%	Acceptable					200	-100%	Fails		(±15)	45
1,3-Dinitrobenzene	19.18	18240	207.1000<	200	4%	Acceptable					200	-100%	Fails		(±15)	45
TETRYL	40.38	18246	209.7000<	200	5%	Acceptable					200	-100%	Fails		(±15)	45
Nitrobenzene	16.47	8291	209.4000<	200	5%	Acceptable					200	-100%	Fails		(±15)	45
2,4,6-Trinitrotoluene	32.59	16243	200.3000<	200	0%	Acceptable					200	-100%	Fails		(±15)	45
4-AM-2,6-DNT	29.08	13363	199.6000<	200	0%	Acceptable					200	-100%	Fails		(±15)	45
2-AM-4,6-DNT	28.74	14987	201.7000<	200	1%	Acceptable					200	-100%	Fails		(±15)	45
2,6-Dinitrotoluene	26.81	10167	203.1000<	200	2%	Acceptable					200	-100%	Fails		(±15)	45
2,4-Dinitrotoluene	25.99	16528	208.1000<	200	4%	Acceptable					200	-100%	Fails		(±15)	45
2-Nitrotoluene	22.38	10164	421.3000<	400	5%	Acceptable					400	-100%	Fails		(±15)	45
<b>4-Nitrotoluene</b>				<b>400</b>	<b>-100%</b>	<b>Fails</b>					<b>400</b>	<b>-100%</b>	<b>Fails</b>		<b>(±15)</b>	
3-Nitrotoluene	22.90	6323	201.4000<	200	1%	Acceptable					200	-100%	Fails		(±15)	45
<b>Nitroglycerin</b>				<b>200</b>	<b>-100%</b>	<b>Fails</b>		38.25	11803	209.9000<	200	5%	Acceptable		(±15)	45
<b>PETN</b>				<b>200</b>	<b>-100%</b>	<b>Fails</b>		48.78	21503	208.1000<	200	4%	Acceptable		(±15)	45
3,5-Dinitroaniline	24.65	13757	205.0000<	200	3%	Acceptable					200	-100%	Fails		(±15)	45
<b>EGDN</b>				<b>200</b>	<b>-100%</b>	<b>Fails</b>		17.80	9381	221.8000	200	11%	Acceptable		(±15)	45

Notes: M = Manually Integrated Signals Differ by More Than 40%  
 D = Operator Disabled Result Signals Differ by More Than 50%  
 O = Over Calibration Range

Chromatography Summary

Injection Date: 7/11/2009 16.16 Operator: fhk  
 DataFile: LC9 I\07102009A BAC-000023.D Vial Num: 4  
 Instrument ID: LC9

Method 8330 Target Analyte Results

Sample: **STD\_5 09GCSV0238 .2K/1/1/1/1**

Method File: LC9.I\07102009A B\8330CNAB.M  
 Start Cal Date: 6/9/2009 20:53 End Cal Date: 6/10/2009 4:20

Matrix: NONE SubList: CAL.sub SpikeList:  
 Samp. Info: STD\_5 09GCSV0238 .2K/1/1/1/1,2  
 Misc. Info: ;5;;;3,CAL.sub;;0,1,0

Dilution Factor	Extract Volume	Sample Volume	Sample Weight
1X	0 mL	0 mL	0 g

Signal 1 UV 250-265								Signal 2 UV 358-205								
Compound Name	RT	Response	PPB	Spike Level	%D	Result	Flag	RT	Response	PPB	Spike Level	%D	Result	Flag	Limits(%)	Flag
3,4-Dinitrotoluene	30.65	4159	104.9000<	100	5%	Acceptable					100	-100%	Fails		(±15)	45
HMX	34.66	3754	91.1000<	100	-9%	Acceptable					100	-100%	Fails		(±15)	45
RDX	25.13	4283	100.6000<	100	1%	Acceptable					100	-100%	Fails		(±15)	45
<b>Picric ACID</b>				<b>200</b>	<b>-100%</b>	<b>Fails</b>					<b>200</b>	<b>-100%</b>	<b>Fails</b>		<b>(±15)</b>	
1,3,5-Trinitrobenzene	21.50	6847	102.2000<	100	2%	Acceptable					100	-100%	Fails		(±15)	45
1,3-Dinitrobenzene	19.16	8905	101.1000<	100	1%	Acceptable					100	-100%	Fails		(±15)	45
TETRYL	40.39	8996	103.4000<	100	3%	Acceptable					100	-100%	Fails		(±15)	45
Nitrobenzene	16.46	4129	104.3000<	100	4%	Acceptable					100	-100%	Fails		(±15)	45
2,4,6-Trinitrotoluene	32.52	7906	97.7300<	100	-2%	Acceptable					100	-100%	Fails		(±15)	45
4-AM-2,6-DNT	29.04	6374	102.7000<	100	3%	Acceptable					100	-100%	Fails		(±15)	45
2-AM-4,6-DNT	28.71	7546	101.6000<	100	2%	Acceptable					100	-100%	Fails		(±15)	45
2,6-Dinitrotoluene	26.77	5040	100.7000<	100	1%	Acceptable					100	-100%	Fails		(±15)	45
2,4-Dinitrotoluene	25.95	8071	101.6000<	100	2%	Acceptable					100	-100%	Fails		(±15)	45
2-Nitrotoluene	22.34	5089	210.9000<	200	5%	Acceptable					200	-100%	Fails		(±15)	45
<b>4-Nitrotoluene</b>				<b>200</b>	<b>-100%</b>	<b>Fails</b>					<b>200</b>	<b>-100%</b>	<b>Fails</b>		<b>(±15)</b>	
3-Nitrotoluene	22.86	3157	100.5000<	100	1%	Acceptable					100	-100%	Fails		(±15)	45
<b>Nitroglycerin</b>				<b>100</b>	<b>-100%</b>	<b>Fails</b>		38.25	5728	101.8000<	100	2%	Acceptable		(±15)	45
<b>PETN</b>				<b>100</b>	<b>-100%</b>	<b>Fails</b>		48.34	10890	105.4000<	100	5%	Acceptable		(±15)	45
3,5-Dinitroaniline	24.62	6884	102.6000<	100	3%	Acceptable					100	-100%	Fails		(±15)	45
<b>EGDN</b>				<b>100</b>	<b>-100%</b>	<b>Fails</b>		17.78	4429	104.7000	100	5%	Acceptable		(±15)	45

Notes M = Manually Integrated Signals Differ by More Than 40%  
 D = Operator Disabled Result Signals Differ by More Than 50%  
 O = Over Calibration Range

TestAmerica West Sacramento

A-000015.D

Chromatography Summary

Injection Date: 7/3/2009 3:17 Operator: fbk  
 DataFile: LC10\07022009\BA-000015.D Vial Num: 21  
 Instrument ID: LC10

Method 8330 Target Analyte Results

Sample: LFV9A1AC 9182192 G9F300242-1 1X

Method File: LC10\07022009\B\8330AB.M  
 Start Cal Date: 6/10/2009 15:33 End Cal Date: 6/11/2009 13:47

Matrix: WATER SubList: WATER sub SpikeList:  
 Samp. Info: LFV9A1AC 9182192 G9F300242-1 IX,0,  
 Misc. Info: ;,1014 4,;20;1;WATER sub,,0,1,LFV9A1AC

Dilution Factor	Extract Volume	Sample Volume	Sample Weight
1X	20 mL	1014.4 mL	0 g

Signal 1 UV 250-265						Signal 2 UV 358-205							
Compound Name	RT	Diff	Response	Conc (ug/L)	Flag	RT	Diff	Response	Conc (ug/L)	Flag	MDL	RL	Flag
3,4-Dinitrotoluene	18.26	0.044	6110	2.5540<		18.26	0.017	11761	2.6820		0.0000	0.00	
HMX											0.0266	0.10	
RDX											0.0641	0.10	
Picric ACID											0.1232	0.99	
1,3,5-Trinitrobenzene	10.11	-0.106	242	0.0323<							0.0306	0.10	45
1,3-Dinitrobenzene											0.0493	0.10	
PIETRYL											0.0493	0.10	
Nitrobenzene											0.0493	0.10	
2,4,6-Trinitrotoluene											0.0237	0.10	
4-AM-2,6-DNT											0.0217	0.10	
2-AM-4,6-DNT											0.0986	0.20	
2,6-Dinitrotoluene											0.0493	0.10	
2,4-Dinitrotoluene											0.0493	0.10	
2-Nitrotoluene											0.0710	0.49	
4-Nitrotoluene											0.0710	0.49	
3-Nitrotoluene											0.0611	0.49	
Nitroglycerin											0.3253	0.64	
PETN											0.2957	0.64	
3,5-Dinitroaniline											0.0246	0.99	

NO

Surrogates:	Spiked	Recovered	% Rec	Spiked	Recovered	% Rec	Limits
3,4-Dinitrotoluene	2.4645	2.5540	104	2.4645	2.6820	109	(79-111)

Notes: M = Manually Integrated Signals Differ by More Than 40%  
 D = Operator Disabled Result Signals Differ by More Than 50%  
 O = Over Calibration Range

Chromatography Summary

Injection Date: 7/3/2009 4:07 Operator: fbk  
Data File: LC10.107022009.BVA-000016.D Vial Num: 22  
Instrument ID: LC10

Method 8330 Target Analyte Results

Sample: LFV9G1AC 9182192 G9F300242-2 1X

Method File: LC10.107022009.BA8330AB.M  
Start Cal Date: 6/10/2009 15:33 End Cal Date: 6/11/2009 13:47

Matrix: WATER SubList: WATER sub SpikeList:  
Samp. Info: 1FV9G1AC 9182192 G9F300242-2 1X,0.  
Misc. Info: ;;975.28;;20;1,WATER sub;0,1,LFV9G1AC

Dilution Factor	Extract Volume	Sample Volume	Sample Weight
1X	20 mL	975.28 mL	0 g

Signal 1 UV 250-265						Signal 2 UV 358-205							
Compound Name	RT	Diff	Response	Conc (ug/L)	Flag	RT	Diff	Response	Conc (ug/L)	Flag	MDL	RL	Flag
3,4-Dinitrotoluene	18.24	0.023	6196	2.6940<		18.24	-0.005	11933	2.8310		0.0000	0.00	
HMX											0.0277	0.10	
RDX											0.0666	0.10	
Picric ACID											0.1282	1.03	
1,3,5-Trinitrobenzene	10.14	-0.074	250	0.0347<							0.0318	0.10	45
1,3-Dinitrobenzene											0.0513	0.10	
TETRYL											0.0513	0.10	
Nitrobenzene											0.0513	0.10	
2,4,6-Trinitrotoluene											0.0246	0.10	
4-AM-2,6-DNT											0.0226	0.10	
2-AM-4,6-DNT											0.1025	0.21	
2,6-Dinitrotoluene											0.0513	0.10	
2,4-Dinitrotoluene											0.0513	0.10	
2-Nitrotoluene											0.0738	0.51	
4-Nitrotoluene											0.0738	0.51	
3-Nitrotoluene											0.0636	0.51	
Nitroglycerin											0.3384	0.67	
PETN											0.3076	0.67	
3,5-Dinitroaniline											0.0256	1.03	

ND

Surrogates	Spiked	Recovered	% Rec	Spiked	Recovered	% Rec	Limits
3,4-Dinitrotoluene	2.5634	2.6940	105	2.5634	2.8310	110	(79-111)

Notes: M = Manually Integrated Signals Differ by More Than 40%  
D = Operator Disabled Result Signals Differ by More Than 50%  
O = Over Calibration Range

Chromatography Summary

Injection Date: 7/3/2009 4:58 Operator: fhk  
Data File: LC10 1107022009.BVA-000017.D Vial Num: 23  
Instrument ID: LC10

Method 8330 Target Analyte Results

Sample: Lfv9J1AC 9182192 G9F300242-3 1X

Method File: LC10 1107022009.BVA330AB.M  
Start Cal Date: 6/10/2009 15:33 End Cal Date: 6/11/2009 13:47

Matrix: WATER SubList: WATER sub SpikeList:  
Samp. Info: Lfv9J1AC 9182192 G9F300242-3 1X,0,  
Misc. Info: ,,1018 21;;20;1;WATER.sub;;0.1;Lfv9J1AC

Dilution Factor	Extract Volume	Sample Volume	Sample Weight
1X	20 mL	1018.21 mL	0 g

Signal 1 UV 250-265						Signal 2 UV 358-205							
Compound Name	RT	Diff	Response	Conc (ug/L)	Flag	RT	Diff	Response	Conc (ug/L)	Flag	MDL	RL	Flag
3,4-Dinitrotoluene	18.25	0.036	6078	2.5310<		18.25	0.009	11702	2.6590		0.0000	0.00	
HMX											0.0265	0.10	
RDX											0.0638	0.10	
Picric ACID											0.1228	0.98	
1,3,5-Trinitrobenzene	10.13	-0.080	276	0.0367<							0.0304	0.10	45
1,3-Dinitrobenzene											0.0491	0.10	
TETRYL											0.0491	0.10	
Nitrobenzene											0.0491	0.10	
2,4,6-Trinitrotoluene											0.0236	0.10	
4-AM-2,6-DNT											0.0216	0.10	
2-AM-4,6-DNT											0.0982	0.20	
2,6-Dinitrotoluene											0.0491	0.10	
2,4-Dinitrotoluene											0.0491	0.10	
2-Nitrotoluene											0.0707	0.49	
4-Nitrotoluene											0.0707	0.49	
3-Nitrotoluene											0.0609	0.49	
Nitroglycerin											0.3241	0.64	
PETN											0.2946	0.64	
3,5-Dinitroaniline											0.0246	0.98	

ND

Surrogates:	Spiked	Recovered	% Rec	Spiked	Recovered	% Rec	Limits
3,4-Dinitrotoluene	2.4553	2.5310	103	2.4553	2.6590	108	(79-111)

Notes: M = Manually Integrated Signals Differ by More Than 40%  
D = Operator Disabled Result Signals Differ by More Than 50%  
O = Over Calibration Range

23LF4 MW7B

A-000018.D

Chromatography Summary

Injection Date: 7/3/2009 5:49 Operator: fmk  
 Data File: LC10.N07022009.B\A-000018.D Vial Num: 24  
 Instrument ID: LC10

Method 8330 Target Analyte Results

Method File: LC10.N07022009.B\8330AB.M  
 Start Cal Date: 6/10/2009 15:33 End Cal Date: 6/11/2009 13:47

Sample: Lfv9K1AC 9182192 G9F300242-4 1X

Dilution Factor	Extract Volume	Sample Volume	Sample Weight
1X	20 mL	1028.11 mL	0 g

Matrix: WATER SubList: WATER sub SpikeList:  
 Samp. Info: 1 FV9K1AC 9182192 G9F300242-4 1X,0,  
 Misc. Info: ;,1028 11,,20,1;WATER sub:,0,1;Lfv9K1AC

Signal 1 UV 250-265						Signal 2 UV 358-205							
Compound Name	RT	Diff	Response	Conc (ug/L)	Flag	RT	Diff	Response	Conc (ug/L)	Flag	MDL	RL	Flag
3,4-Dinitrotoluene	18.24	0.028	5788	2.3870<		18.24	0.001	11100	2.4980		0.0000	0.00	
HMX											0.0263	0.10	
RDX											0.0632	0.10	
Picric ACID											0.1216	0.97	
1,3,5-Trinitrobenzene	10.17	-0.045	264	0.0348<							0.0302	0.10	45
1,3-Dinitrobenzene											0.0486	0.10	
TETRYL											0.0486	0.10	
Nitrobenzene											0.0486	0.10	
2,4,6-Trinitrotoluene											0.0233	0.10	
4-AM-2,6-DNT											0.0214	0.10	
2-AM-4,6-DNT											0.0973	0.19	
2,6-Dinitrotoluene											0.0486	0.10	
2,4-Dinitrotoluene											0.0486	0.10	
2-Nitrotoluene											0.0700	0.49	
4-Nitrotoluene											0.0700	0.49	
3-Nitrotoluene											0.0603	0.49	
Nitroglycerin											0.3210	0.63	
PETN											0.2918	0.63	
3,5-Dinitroaniline											0.0243	0.97	

NO

Surrogates:	Spiked	Recovered	% Rec	Spiked	Recovered	% Rec	Limits
3,4-Dinitrotoluene	2.4316	2.3870	98	2.4316	2.4980	103	(79-111)

Notes: M = Manually Integrated Signals Differ by More Than 40%  
 D = Operator Disabled Result Signals Differ by More Than 50%  
 O = Over Calibration Range



23LF4MWSA

TestAmerica West Sacramento

A-000021.D

Chromatography Summary

Injection Date: 7/3/2009 8:20 Operator: fhk
Data File: LC10 I07022009 BVA-000021.D Vial Num: 27
Instrument ID: LC10

Method 8330 Target Analyte Results

Sample: Lfv9M1AC 9182192 G9F300242-5 1X Method File: LC10 I07022009 B8330AB.M
Start Cal Date: 6/10/2009 15:33 End Cal Date: 6/11/2009 13:47

Matrix: WATER SubList: WATER sub SpikeList:
Samp. Info: Lfv9M1AC 9182192 G9F300242-5 1X,0;
Misc. Info: .:1016 72,,20:1;WATER sub..0.1.Lfv9M1AC

Table with 4 columns: Dilution Factor, Extract Volume, Sample Volume, Sample Weight. Values: 1X, 20 mL, 1016.72 mL, 0 g

Main chromatography table with columns: Compound Name, RT, Diff, Response, Conc (ug/L), Flag, RT, Diff, Response, Conc (ug/L), Flag, MDL, RI, Flag. Includes handwritten notes like 'converted on 6/21/09'.

Surrogate table with columns: Surrogate, Spiked, Recovered, % Rec, Spiked, Recovered, % Rec, Limits. Row: 3,4-Dinitrotoluene, 2.4589, 2.5880, 105, 2.4589, 2.7120, 110, (79-111)

Notes: M = Manually Integrated Signals Differ by More Than 40%
D = Operator Disabled Result Signals Differ by More Than 50%
O = Over Calibration Range

Chromatography Summary

Injection Date: 7/11/2009 15:14 Operator: fmk  
 DataFile: LC9 I07102009A.BVC-000022.D Vial Num: 28  
 Instrument ID: LC9

Method 8330 Target Analyte Results

Sample: LFV9M1AC 9182192 G9F300242-5 1X Method File: LC9 I07102009A.BW330CNAB.M  
 Start Cal Date: 6/9/2009 20:53 End Cal Date: 6/10/2009 4:20

Matrix: WATER SubList: WATER sub SpikeList:  
 Samp. Info: LFV9M1AC 9182192 G9F300242-5 1X;0;  
 Misc. Info: ;,1016.72;20;1,WATER sub;0;1

Dilution Factor	Extract Volume	Sample Volume	Sample Weight
1X	20 mL	1016.72 mL	0 g

Signal 1 UV 250-265						Signal 2 UV 358-205							
Compound Name	RT	Diff	Response	Conc (ug/L)	Flag	RT	Diff	Response	Conc (ug/L)	Flag	MDL	RL	Flag
3,4-Dinitrotoluene	30.65	-0.044	5665	2.8100<							0.0000	0.00	45
HMX	34.67	-0.007	430	0.2052<							0.0266	0.10	45
RDX	25.15	-0.016	8165	3.7740<							0.0639	0.10	45
Picric ACID											0.1229	0.97	
1,3,5-Trinitrobenzene											0.0305	0.10	
1,3-Dinitrobenzene											0.0492	0.10	
TETRYL											0.0492	0.10	
Nitrobenzene											0.0492	0.10	
2,4,6-Trinitrotoluene											0.0236	0.10	
4-AM-2,6-DNT											0.0216	0.10	
2-AM-4,6-DNT											0.0984	0.19	
2,6-Dinitrotoluene											0.0492	0.10	
2,4-Dinitrotoluene											0.0492	0.10	
2-Nitrotoluene											0.0708	0.48	
4-Nitrotoluene											0.0708	0.48	
3-Nitrotoluene											0.0610	0.48	
Nitroglycerin											0.3246	0.63	
PETN											0.2951	0.63	
3,5-Dinitroaniline											0.0246	0.97	

*continued*  
*7/11/09*

Surrogates:	Spiked	Recovered	% Rec	Spiked	Recovered	% Rec	Limits
3,4-Dinitrotoluene	2.4589	2.8100	114	2.4589	0	0	(48-143)

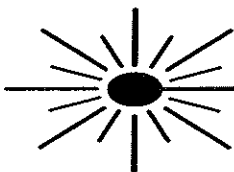
Notes: M = Manually Integrated Signals Differ by More Than 40%  
 D = Operator Disabled Result Signals Differ by More Than 50%  
 O = Over Calibration Range

**Data Validation Report**

**Michael Baker Jr., Inc**

**Camp Bonneville**

**SDG#: PSF0932**

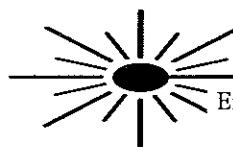


**ECT.CON INC.**

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Environmental and Computer  
Technology Consultants

3531 Fox Chase Drive  
Imperial, PA 15126  
(724) 695-8042  
FAX (724) 695-2698  
e-mail: [ectconinc@comcast.net](mailto:ectconinc@comcast.net)



## Data Validation Report

SDG#	PSF0932
Validation Report Date	August 17, 2009
Validation Guidance	USEPA CLP National Functional Guidelines for Data Review
Client Name	Michael Baker Jr., Inc.
Project Name	Camp Bonneville
Laboratory	TestAmerica
Analytical Parameters	Explosives
Analytical Method	SW-846 8330

### Samples/Matrix:

Date Sampled	Location	Portland ID	Sacramento ID	Explosives	Matrix
06/29/09	23L4MW01AW	PSF0932-01	LFXXW	X	Aqueous
06/29/09	23L4MW01BW	PSF0932-02	LFXX6	X	Aqueous
06/26/09	23L4MW460W	PSF0932-04	LFX11	X	Aqueous
06/26/09	23L4MW03BW	PSF0932-05	LFX12	X	Aqueous
06/26/09	23L4MW03AW	PSF0932-06	LFX15	X	Aqueous

Analytical data in this report were screened to determine analytical limitations of the data based on specific quality control criteria. This screening assumes analytical results are correct as reported and merely provides an interpretation of the reported quality control results. Laboratory calculations have been verified as part of this validation. Specific findings on analytical limitations are presented in this report. Annotated Form 1s or spreadsheets for samples reviewed are included after the Data Assessment Findings. Form 1s for the MS/MSD samples and spreadsheets are not annotated.

### SUMMARY

The sample set for the Camp Bonneville site consists of five aqueous field samples. These samples were analyzed for explosives as provided in the Table 1. The findings presented in this review of the analytical data assume that the information presented by the analytical laboratory is correct.

The explosives findings are based upon the assessment of the following:

- \* ● Data Completeness
- \* ● Holding Times
- \* ● Calibration (Initial and Continuing)
- Blanks
- \* ● System Monitoring Compounds (Surrogate Spikes)
- \* ● Matrix Spike/Matrix Spike Duplicates
- \* ● Laboratory Control Standard Results
- \* ● Target Compound Identification
- Compound Quantification and Reported Contract Quantitation Limits

- \* • System Performance
- \* Criteria were met for this evaluation item.

This evaluation was conducted in accordance with USEPA CLP National Functional Guidelines for Organic Data Review and the analytical method. Findings from this evaluation should be considered when using the analytical data. This report presents a summary of the data qualifications based on the review of the aforementioned evaluation criteria. This is followed by annotated Form 1s/ spreadsheets. Finally, the worksheets used to perform the evaluation are provided.

## **FINDINGS**

### **1. Blanks**

The laboratory method blank exhibited contamination for the following parameter:

Blank	Compound	Maximum Concentration (ppb)	Action Limit (ppm)	Action
LF1EN1	1,3,5-Trinitrobenzene	0.2088	1.044	U sample results < RL

RL - reporting limit

### **2. Compound Quantitation**

Positive results less than the reporting limit were qualified as estimated "J" due to the level of uncertainty.

In sample 23L4MW03BW, the positive result below the reporting limit for HMX was not confirmed. Therefore, the form one was amended to show the result as nondetected "U" at the reporting limit.

## **NOTES**

### **Matrix Spike/Matrix Spike Duplicate Results**

A MS/MSD was not analyzed with this SDG. A laboratory control sample was analyzed. Data were not qualified on this basis.

### **Field Duplicate Results**

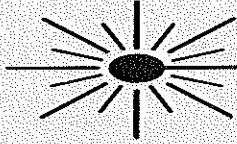
No field duplicates were included with this SDG. Data were not qualified on this basis.

  
Data Reviewer

8/12/09  
Date

### **Glossary of Data Qualifiers**

U	Not Detected.	The associated number indicates approximate sample concentration necessary to be detected.
UJ	Not Detected.	Quantitation limit may be inaccurate or imprecise.
J	Analyte Present.	Reported value may not be accurate or precise.
N	Consider Present.	Tentative identification. Special methods may be needed to confirm its presence or absence in future sampling efforts.
R	Unusable Result.	Analyte may or may not be present in the sample.
UR	Unusable Result.	Analyte may or may not be present in the sample.



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*Annotated Form 1's*  
*(Spreadsheet)*

TestAmerica Portland

Client Sample ID: PSF0932-01

HPLC *2324MW01AW*

Lot-Sample #...: G9G010225-001    Work Order #...: LFXXW1AC    Matrix.....: WATER  
 Date Sampled...: 06/29/09    Date Received...: 07/01/09  
 Prep Date.....: 07/02/09    Analysis Date...: 07/08/09  
 Prep Batch #...: 9183251  
 Dilution Factor: 0.97    Method.....: SW846 8330

PARAMETER	RESULT	REPORTING LIMIT	UNITS	MDL
Nitroglycerin	ND	0.63	ug/L	0.15
PETN	ND	0.63	ug/L	0.22
Picric Acid	ND	0.97	ug/L	0.12
2-Amino-4,6-dinitrotoluene	ND	0.19	ug/L	0.097
4-Amino-2,6-dinitrotoluene	ND	0.097	ug/L	0.021
1,3-Dinitrobenzene	ND	0.097	ug/L	0.048
2,4-Dinitrotoluene	ND	0.097	ug/L	0.048
2,6-Dinitrotoluene	ND	0.097	ug/L	0.048
HMX	ND	0.097	ug/L	0.026
Nitrobenzene	ND	0.097	ug/L	0.048
2-Nitrotoluene	ND	0.39	ug/L	0.070
3-Nitrotoluene	ND	0.39	ug/L	0.060
4-Nitrotoluene	ND	0.48	ug/L	0.070
RDX	0.13	0.097	ug/L	0.063
Tetryl	ND	0.097	ug/L	0.048
1,3,5-Trinitrobenzene	<del>0.033</del>	0.097 U	ug/L	0.030
2,4,6-Trinitrotoluene	ND	0.097	ug/L	0.023

SURROGATE	PERCENT RECOVERY	RECOVERY LIMITS
3,4-Dinitrotoluene	91	(79 - 111)

NOTE(S) :

1 Estimated result Result is less than RL.

*UK  
8/10/09*



TestAmerica Portland

Client Sample ID: PSF0932-02

HPLC *23LYMWO18W*

Lot-Sample #...: G9G010225-002    Work Order #...: LFXX61AC    Matrix.....: WATER  
 Date Sampled...: 06/29/09    Date Received...: 07/01/09  
 Prep Date.....: 07/02/09    Analysis Date...: 07/08/09  
 Prep Batch #...: 9183251  
 Dilution Factor: 0.97    Method.....: SW846 8330

PARAMETER	RESULT	REPORTING		
		LIMIT	UNITS	MDL
Nitroglycerin	ND	0.63	ug/L	0.15
PETN	ND	0.63	ug/L	0.22
Picric Acid	ND	0.97	ug/L	0.12
2-Amino-4,6-dinitrotoluene	ND	0.19	ug/L	0.097
4-Amino-2,6-dinitrotoluene	ND	0.097	ug/L	0.021
1,3-Dinitrobenzene	ND	0.097	ug/L	0.048
2,4-Dinitrotoluene	ND	0.097	ug/L	0.048
2,6-Dinitrotoluene	ND	0.097	ug/L	0.048
HMX	ND	0.097	ug/L	0.026
Nitrobenzene	ND	0.097	ug/L	0.048
2-Nitrotoluene	ND	0.39	ug/L	0.070
3-Nitrotoluene	ND	0.39	ug/L	0.060
4-Nitrotoluene	ND	0.48	ug/L	0.070
RDX	ND	0.097	ug/L	0.063
Tetryl	ND	0.097	ug/L	0.048
1,3,5-Trinitrobenzene	<del>0.033</del> <i>3</i>	0.097 <i>U</i>	ug/L	0.030
2,4,6-Trinitrotoluene	ND	0.097	ug/L	0.023
	PERCENT	RECOVERY		
SURROGATE	RECOVERY	LIMITS		
3,4-Dinitrotoluene	93	(79 - 111)		

NOTE(S):

*I* Estimated result. Result is less than RL.

*UK  
8/16/09*

TestAmerica Portland

Client Sample ID: PSF0932-04

HPLC 234 MW 960W

Lot-Sample #...: G9G010225-003 Work Order #...: LFX111AC Matrix.....: WATER  
 Date Sampled...: 06/26/09 Date Received...: 07/01/09  
 Prep Date.....: 07/02/09 Analysis Date...: 07/08/09  
 Prep Batch #...: 9183251  
 Dilution Factor: 0.97 Method.....: SW846 8330

PARAMETER	RESULT	REPORTING		
		LIMIT	UNITS	MDL
Nitroglycerin	ND	0.63	ug/L	0.15
PETN	ND	0.63	ug/L	0.22
Picric Acid	ND	0.97	ug/L	0.12
2-Amino-4,6-dinitrotoluene	ND	0.19	ug/L	0.097
4-Amino-2,6-dinitrotoluene	ND	0.097	ug/L	0.021
1,3-Dinitrobenzene	ND	0.097	ug/L	0.048
2,4-Dinitrotoluene	ND	0.097	ug/L	0.048
2,6-Dinitrotoluene	ND	0.097	ug/L	0.048
HMX	0.029 J J	0.097	ug/L	0.026
Nitrobenzene	ND	0.097	ug/L	0.048
2-Nitrotoluene	ND	0.39	ug/L	0.070
3-Nitrotoluene	ND	0.39	ug/L	0.060
4-Nitrotoluene	ND	0.48	ug/L	0.070
RDX	4.1	0.097	ug/L	0.063
Tetryl	ND	0.097	ug/L	0.048
1,3,5-Trinitrobenzene	ND	0.097	ug/L	0.030
2,4,6-Trinitrotoluene	ND	0.097	ug/L	0.023

SURROGATE	PERCENT	RECOVERY
	RECOVERY	LIMITS
3,4-Dinitrotoluene	93	(79 - 111)

NOTE(S):  
 J Estimated result. Result is less than RL.

WJ  
8/6/09

TestAmerica Portland

Client Sample ID: PSF0932-05

HPLC

23LY NW03BW

Lot-Sample #...: G9G010225-004    Work Order #...: LFX121AC    Matrix.....: WATER  
 Date Sampled...: 06/26/09    Date Received...: 07/01/09  
 Prep Date.....: 07/02/09    Analysis Date...: 07/08/09  
 Prep Batch #...: 9183251  
 Dilution Factor: 0.98    Method.....: SW846 8330

PARAMETER	RESULT	REPORTING		
		LIMIT	UNITS	MDL
Nitroglycerin	ND	0.64	ug/L	0.15
PETN	ND	0.64	ug/L	0.23
Picric Acid	ND	0.98	ug/L	0.12
2-Amino-4,6-dinitrotoluene	ND	0.20	ug/L	0.098
4-Amino-2,6-dinitrotoluene	ND	0.098	ug/L	0.022
1,3-Dinitrobenzene	ND	0.098	ug/L	0.049
2,4-Dinitrotoluene	ND	0.098	ug/L	0.049
2,6-Dinitrotoluene	ND	0.098	ug/L	0.049
HMX	<del>0.027</del>	0.098	ug/L	0.026
Nitrobenzene	ND	0.098	ug/L	0.049
2-Nitrotoluene	ND	0.39	ug/L	0.071
3-Nitrotoluene	ND	0.39	ug/L	0.061
4-Nitrotoluene	ND	0.49	ug/L	0.071
RDX	4.1	0.098	ug/L	0.064
Tetryl	ND	0.098	ug/L	0.049
1,3,5-Trinitrobenzene	ND	0.098	ug/L	0.030
2,4,6-Trinitrotoluene	ND	0.098	ug/L	0.024
	PERCENT	RECOVERY		
<u>SURROGATE</u>	<u>RECOVERY</u>	<u>LIMITS</u>		
3,4-Dinitrotoluene	93	(79 - 111)		

NOTE(S) :

J Estimated result Result is less than RL.

UK  
8/10/09

TestAmerica Portland

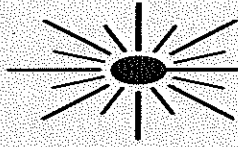
Client Sample ID: PSF0932-06

HPLC *234MW03AW*

Lot-Sample #...: G9G010225-005    Work Order #...: LFX151AC    Matrix.....: WATER  
 Date Sampled...: 06/26/09    Date Received...: 07/01/09  
 Prep Date.....: 07/02/09    Analysis Date...: 07/08/09  
 Prep Batch #...: 9183251  
 Dilution Factor: 0.97    Method.....: SW846 8330

PARAMETER	RESULT	REPORTING LIMIT	UNITS	MDL
Nitroglycerin	ND	0.63	ug/L	0.15
PETN	ND	0.63	ug/L	0.22
Picric Acid	ND	0.97	ug/L	0.12
2-Amino-4,6-dinitrotoluene	ND	0.19	ug/L	0.097
4-Amino-2,6-dinitrotoluene	ND	0.097	ug/L	0.021
1,3-Dinitrobenzene	ND	0.097	ug/L	0.048
2,4-Dinitrotoluene	ND	0.097	ug/L	0.048
2,6-Dinitrotoluene	ND	0.097	ug/L	0.048
HMX	0.38	0.097	ug/L	0.026
Nitrobenzene	ND	0.097	ug/L	0.048
2-Nitrotoluene	ND	0.39	ug/L	0.070
3-Nitrotoluene	ND	0.39	ug/L	0.060
4-Nitrotoluene	ND	0.48	ug/L	0.070
RDX	9.4	0.097	ug/L	0.063
Tetryl	ND	0.097	ug/L	0.048
1,3,5-Trinitrobenzene	ND	0.097	ug/L	0.030
2,4,6-Trinitrotoluene	ND	0.097	ug/L	0.023
	PERCENT	RECOVERY		
SURROGATE	RECOVERY	LIMITS		
3,4-Dinitrotoluene	95	(79 - 111)		

*WST  
6/10/09*



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## *Support Documentation*

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## *Laboratory Case Narrative*

## **DATA DELIVERABLES PACKAGE**

Michael Baker Jr., Inc.  
James D. Peyton  
5261 Fountain Drive, Suite A  
Crown Point, IN 46307

Client Project: Camp Bonneville Groundwater  
Client Project Number: 110002 Task 6200

TA Work Order#: PSF0932  
TA Project Manager: Estella Rieben

The total number of pages contained in this data package is:

163

July 31, 2009

TestAmerica - Portland  
9405 S. W. Nimbus Avenue  
Beaverton, Oregon 97008  
(503) 906-9200  
(503) 906-9210

## CASE NARRATIVE

**Client:** Michael Baker Jr., Inc. **Date Sampled:** 6/26/2009 and 6/29/2009  
**Project:** Camp Bonneville Groundwater 110002 Task 6200 **Date Received:** 6/29/2009  
**Lab:** PSF0932

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**SAMPLE RECEIPT:** Samples were received intact, on ice, with chain of custody documentation. The sample temperature was measured at 2.3° C upon receipt at the laboratory.

**HOLDING TIMES:** All samples were analyzed within holding times.

**PROBLEMS ENCOUNTERED:** No problems were encountered.

**QA/QC CRITERIA:** No problems were encountered during sample analysis.

**OBSERVATIONS:** No significant observations were made.

**SUBCONTRACTED:** 314.0\_9196313 (Perchlorate), 314.0\_9196314 (Perchlorate), and SW846 8330 analyses were performed at the TestAmerica facility in West Sacramento, CA. The data package for the specified analyses is included in this data deliverables package as Appendix A.

*"I certify that this data package is in compliance with the contract both technically and for completeness, for all conditions other than the conditions detailed above. Release of the data contained in this data package has been authorized by the Laboratory Director or his designee, as verified by the following signature."*

**TestAmerica Portland**



Estella Rieben  
Project Manager



**Michael Baker Jr., Inc.**  
5261 Fountain Drive, Suite A  
Crown Point, IN 46307

Project Name: **Camp Bonneville Groundwater**  
Project Number: 110002 Task 6200  
Project Manager: James D. Peyton

Report Created:  
07/21/09 08:53

## ANALYTICAL REPORT FOR SAMPLES

Sample ID	Laboratory ID	Matrix	Date Sampled	Date Received
23L4MW01AW	PSF0932-01	Water	06/29/09 10:35	06/29/09 17:00
23L4MW01BW	PSF0932-02	Water	06/29/09 14:40	06/29/09 17:00
TB-261	PSF0932-03	Water	06/29/09 00:00	06/29/09 17:00
23L4MW460W	PSF0932-04	Water	06/26/09 15:00	06/29/09 17:00
23L4MW03BW	PSF0932-05	Water	06/26/09 17:15	06/29/09 17:00
23L4MW03AW	PSF0932-06	Water	06/26/09 18:15	06/29/09 17:00

TestAmerica Portland



Estella Rieben, Project Manager

*The results in this report apply to the samples analyzed in accordance with the chain of custody document. This analytical report shall not be reproduced except in full, without the written approval of the laboratory.*



# TestAmerica

THE LEADER IN ENVIRONMENTAL TESTING

July 28, 2009

**TestAmerica Project Number: G9G010225**  
PO/Contract: PSF0932

Estella Rieben  
TestAmerica Portland  
Nimbus Corporate Center  
9405 SW Nimbus Ave  
Beaverton, OR 97008

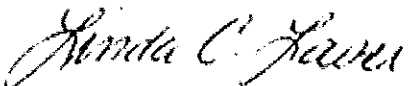
Dear Ms. Rieben,

This report contains the analytical results for the samples received under chain of custody by TestAmerica on July 1, 2009. These samples are associated with your Michael Baker Jr., Inc. project.

The test results in this report meet all NELAC requirements for parameters that accreditation is required or available. Any exceptions to NELAC requirements are noted in the case narrative. The case narrative is an integral part of this report.

If you have any questions, please feel free to call me at (916) 374-4362.

Sincerely,



Linda C. Laver  
Project Manager

## Case Narrative

### TestAmerica West Sacramento Project Number G9G010225

#### General Comments

The chain of custody (coc) indicates that the samples containers were preserved with Hydrochloric Acid (HCL). These analyses do not require acid preservation. The pH of the samples was checked and they were at a pH of 7. According to Ms. Rieben, the coc was incorrect and the samples were not preserved.

#### **WATER, 8330, Nitroaromatics & Nitramines**

Samples: 1, 2, 3, 4, 5

There was insufficient sample volume to prepare a matrix spike/matrix spike duplicate (MS/MSD) pair with this extraction batch.

There are no other anomalies associated with this project.

## Sample Summary

### TestAmerica West Sacramento Project Number G9G010225

<u>WO#</u>	<u>Sample #</u>	<u>Client Sample ID</u>	<u>Sampling Date</u>	<u>Received Date</u>
LFXXW	1	PSF0932-01	6/29/2009 10:35 AM	7/1/2009 09:45 AM
LFXX6	2	PSF0932-02	6/29/2009 02:40 PM	7/1/2009 09:45 AM
LFX11	3	PSF0932-04	6/26/2009 03:00 PM	7/1/2009 09:45 AM
LFX12	4	PSF0932-05	6/26/2009 05:15 PM	7/1/2009 09:45 AM
LFX15	5	PSF0932-06	6/26/2009 08:15 PM	7/1/2009 09:45 AM

#### Notes(s):

- The analytical results of the samples listed above are presented on the following pages.
- All calculations are performed before rounding to avoid round-off errors in calculated results.
- Results noted as "ND" were not detected at or above the stated limit.
- This report must not be reproduced, except in full, without the written approval of the laboratory.
- Results for the following parameters are never reported on a dry weight basis: color, corrosivity, density, flashpoint, ignitability, layers, odor, paint filter test, pH, porosity, pressure, reactivity, redox potential, specific gravity, spot tests, solids, solubility, temperature, viscosity, and weight

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## *Chain of Custody*

# TestAmerica

THE LEADER IN ENVIRONMENTAL TESTING

11720 North Creek Drive N Suite 100, Bothell, WA 98011 8244

19022 E. First Ave, Spokane, WA 99208-5502

9405 NW Nambour Ave, Vancouver, WA 98686-1111

1000 W. Broadway, Suite 100, Portland, OR 97201-3000

125 430-9200 FAX 426-9210  
 300 921-9200 FAX 921-9290  
 300 921-9200 FAX 921-9290

## CHAIN OF CUSTODY REPORT

Work Order #: **PSFO932**

CLIENT: **MICHAEL BAKER JR. INC.**  
 REPORT TO: **MARGARET JAMES**  
 ADDRESS: **5201 FORT WAIN DRIVE, SUITE A**  
**CRAICOR BLDG - IN 46307**  
 PHONE: **219 736 6663** FAX: **219 755 6233**  
 PROJECT NAME:  
 PROJECT NUMBER:  
 SAMPLED BY:

INVOICE TO: **BAKER**  
 ← SAME ADDRESS  
 P.O. NUMBER: **10002** TASK: **6200**  
 PRESERVATIVE:  
 - HCL - **N/A** - **N/A** - **N/A**

CLIENT SAMPLE IDENTIFICATION	SAMPLING DATE/TIME	REQUESTED ANALYSES		OTHER	MATRIX (W, S, O)	# OF CONT.	LOCATION/ COMMENTS	TA VOID
		VOC'S	EXPRESIVES					
1. 23L4 MW01AW	6/29/09 - 10:35	X	X		W	6	NO TICS	
2. 23L4 MW01BW	6/29/09 - 14:40	X	X		W	6	NO TICS	
3. TB-261	6/29/09 - N/A	X	X		W	6	NO TICS	
4. 23L4 MW460W	6/26/09 - 1500	X	X		W	6	NO TICS	
5. 23L4 MW03BW	6/24/09 - 1715	X	X		W	6	NO TICS	
6. 23L4 MW03AW	6/26/09 - 1815	X	X		W	6	NO TICS	
7								
8								
9								
10								

RECEIVED BY: **William Pelkey** DATE: **6/29/09** TIME: **15:15**  
 PRINT NAME: **WILLIAM PELKEY** FIRM: **BAKER**  
 RECEIVED BY: **Bob F** DATE: **6/29/09** TIME: **17:00**  
 PRINT NAME: **BOB F** FIRM: **TAP**  
 RECEIVED BY: **William Pelkey** DATE: **6/29/09** TIME: **17:00**  
 PRINT NAME: **WILLIAM PELKEY** FIRM: **TAP**  
 ADDITIONAL REMARKS:  
 TEMP: **73** OF **73** CHANGE OF

TestAmerica Portland  
**Sample Receiving Checklist**

Work Order #: PSF0932 Date/Time Received: 10/29/09 1700  
 Client Name and Project: Michael Baker

~~G-BA-11E~~ Camp Bonneville GW

Time Zone:  
 EDT/EST     CDT/CST     MDT/MST     PDT/PST     AK     OTHER

**Unpacking Checks:**

Cooler #(s): 1  
 Temperatures: 2.0  
 Digi #1  Digi #2  IR Gun  (  Plastic  Glass )

**Temperature out of Range:**

Not enough or No Ice  
 Ice Melted  
 W/in 4 Hrs of collection  
 Other: \_\_\_\_\_

Initials: MBK

N/A Yes No

- 1. If ESI client, were temp blanks received? If no, document on NOD.
- 2. Cooler Seals intact? (N/A if hand delivered) if no, document on NOD.
- 3. Chain of Custody present? If no, document on NOD.
- 4. Bottles received intact? If no, document on NOD.
- 5. Sample is not multiphasic? If no, document on NOD.
- 6. Proper Container and preservatives used? If no, document on NOD.
- 7. pH of all samples checked and meet requirements? If no, document on NOD.
- 8. Cyanide samples checked for sulfides and meet requirements? If no, notify PM.
- 9. HF Dilution required?
- 10. Sufficient volume provided for all analysis? If no, document on NOD and consult PM before proceeding.
- 11. Did chain of custody agree with samples received? If no, document on NOD.
- 12. Is the "Sampled by" section of the COC completed?
- 13. Were VOA/Oil Syringe samples without headspace?
- 14. Were VOA vials preserved?  HCl  Sodium Thiosulfate  Ascorbic Acid
- 15. Did samples require preservation with sodium thiosulfate?
- 16. If yes to #14, was the residual chlorine test negative? If no, document on NOD.
- 17. Are dissolved/field filtered metals bottles sediment-free? If no, document on NOD.
- 18. Is sufficient volume provided for client requested MS/MSD or matrix duplicates? If no, document on NOD and contact PM before proceeding.
- 19. Are analyses with short holding times received in hold?
- 20. Was Standard Turn Around (TAT) requested?
- 21. Receipt date(s) < 48 hours past the collection date(s)? If no, notify PM.

TestAmerica Portland  
Sample Receiving Checklist

Work Order #: PSF0932

**Login Checks:**

Initials: BUE

N/A Yes No

22. Sufficient volume provided for all analysis? If no, document on NOD & contact PM.
23. Sufficient volume provided for client requested MS/MSD or matrix duplicates? If no, document on NOD and contact PM.
24. Did the chain of custody include "received by" and "relinquished by" signatures, dates and times?
25. Were special log in instructions read and followed?
26. Were tests logged checked against the COC?
27. Were rush notices printed and delivered?
28. Were short hold notices printed and delivered?
29. Were subcontract COCs printed?
30. Was HF dilution logged?

**Labeling and Storage Checks:**

Initials: BUE

N A Yes No

31. Were the subcontracted samples/containers put in Sx fridge?
32. Were sample bottles and COC double checked for dissolved/filtered metals?
33. Did the sample ID, Date, and Time from label match what was logged?
34. Were Foreign sample stickers affixed to each container and containers stored in foreign fridge?
35. Were HF stickers affixed to each container, and containers stored in Sx fridge?
36. Was an NOD for created for noted discrepancies and placed in folder?

Document any problems or discrepancies and the actions taken to resolve them on a Notice of Discrepancy form (NOD).



CLIENT Haz - Portland PM LL LOG # 59283

LOT# (QUANTIMS ID) G9G016225 QUOTE# 83765 LOCATION W89B

DATE RECEIVED 7-1-09 TIME RECEIVED 945 Initials as Date 7-1-09

- DELIVERED BY
- FEDEX
  - AIRBORNE
  - UPS
  - TAL COURIER
  - OTHER
  - CA OVERNIGHT
  - GOLDENSTATE
  - BAX GLOBAL
  - VALLEY LOGISTICS

CUSTODY SEAL STATUS  INTACT  BROKEN  N/A

CUSTODY SEAL #(S) 204345

SHIPPING CONTAINER(S)  TAL  CLIENT  N/A

TEMPERATURE RECORD (IN °C) IR 4  5  OTHER

COC #(S) NA

TEMPERATURE BLANK Observed: 0 Corrected: 2

SAMPLE TEMPERATURE  
Observed: 2 1 2 Average: 2 Corrected Average: 2

COLLECTOR'S NAME  Verified from COC  Not on COC

pH MEASURED  YES  ANOMALY  N/A <sup>on 7-1-09</sup>

LABELED BY .....

LABELS CHECKED BY .....

PEER REVIEW  N/A

SHORT HOLD TEST NOTIFICATION SAMPLE RECEIVING  
WETCHEM  N/A  
VOA-ENCORES  N/A

METALS NOTIFIED OF FILTER/PRESERVE VIA VERBAL & EMAIL  N/A

COMPLETE SHIPMENT RECEIVED IN GOOD CONDITION WITH APPROPRIATE TEMPERATURES, CONTAINERS, PRESERVATIVES  N/A

CLOUSEAU  TEMPERATURE EXCEEDED (2 °C -- 6 °C)  N/A

WET ICE  BLUE ICE  GEL PACK  NO COOLING AGENTS USED  PM NOTIFIED

Notes: PJT & ASB with pH of 7, COC lists HCL

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# *Worksheets*

**HOLDING TIMES**

SAMPLE DATE	SAMPLE ID	Portland LAB ID	Sacramento Lab ID	Matrix	Prep	Anal
06/29/09	23L4MW01AW	PSF0932-01	LFXXW	Aq	07/02/09	7/8/09 1X
06/29/09	23L4MW01BW	PSF0932-02	LFXX6	Aq	07/02/09	7/8/09 1X
06/26/09	23L4MW460W	PSF0932-04	LFX11	Aq	07/02/09	7/8/09 1X
06/26/09	23L4MW03BW	PSF0932-05	LFX12	Aq	07/02/09	7/8/09 1X
06/26/09	23L4MW03AW	PSF0932-06	LFX15	Aq	07/02/09	7/8/09 1X

7 days from collection to extraction  
 Sample Date 6/26/2009 Extract By 7/3/2009  
 7 days from collection to extraction  
 Sample Date 6/29/2009 Extract By 7/6/2009  
 40 days from extraction to analysis  
 Sample Date 7/2/2009 Extract By 8/11/2009

**TARGET COMPOUNDS AND QUANTITATION LIMITS**

Was a target compound list provided by the client? No  
 Did Sample Form 1s match the target compound list NA  
 Were required quantitation limits provided by the client? No  
 Did all compounds meet the required quantitation limits? NA

**SYSTEM MONITORING COMPOUNDS**

SAMPLE	SURROGATE	COLUMN I	DF	ACTION
All IN				

Were surrogate RTs within windows established by the ICAL? Yes  
 Were there any transcription errors between the raw data and Form 2? No  
 Were laboratory acceptance limits used as the basis for validation? 30-150  
 Did the laboratory provide CLP Form II or equivalent? Yes

01AW

SURR	AMOUNT FOUND	AMOUNT SPIKED	% R	F1
34-DNT A	2.233	2.5	89.3	91
34-DNT C	2.456	2.5	98.2	

**MS/MSD RECOVERY and RELATIVE PERCENT DIFFERENCE**

NonePresent

**LABORATORY CONTROL SAMPLES**

SAMPLE	COMPOUND	LCS %R	LCSD %R	RPD	ACTION
LF1EN1	All Parameters	OK	NA	NA	none, ND

Were laboratory acceptance limits used as the basis for validation? Yes  
 Did the laboratory provide CLP Form III or equivalent? Yes  
 Were chromatograms and quan reports present for all LCS/LCDs? Yes

LF1EN1

RDX

All Parameters	AMOUNT FOUND	AMOUNT SPIKED	% R	FORM 3
LCS	0.97	1	97.0	127

%R = (Amount Found/Amount Spiked)\*100

**BLANKS**

BLANK	COMPOUND	RESULT	5X OR 10X	ACTION LEVEL	ACTION
LF1EN1	123-Tnbenzene	0.2088	5	1.044	U

appears on A channel but not on C

NOTE Equipment/Field blanks are not qualified on the basis of laboratory method blank contamination or contamination in other field quality control blanks.

**CALIBRATION**

**Initial**

Are chromatograms and quan reports present for all ICAL standards? Yes  
 Are CLP For VIs or equivalent present and complete? Yes  
 Any transcription or calculation errors? No  
 What linearity criteria was used? 20% RSD or 0.99  
 Were RT window documented? Yes

**Continuing**

Are chromatograms and quan reports present for all CCV standards? Yes  
 Are CLP Form VIIIs or equivalent present and complete? Yes  
 Any transcription or calculation errors? No  
 What %D criteria was used? 25% D  
 Have all samples and standards been listed on an Analytical Sequence? Yes  
 Was a proper analytical sequence followed? Yes

**INITIAL CALIBRATION - A**

Date	6/10/2009	Rptd Avg CF	77.89163	Rptd %RSD	5.156
Instrument	A	Calcd Avg CF	77.892	Calcd %RSD	5.156
Compound	RDX				
Reported CF	80.13	CF1	81.80	STD DEV	4
Calculated CF	80.42	CF2	75.90		
		CF3	77.15		
Response	16084	CF4	80.72		
Conc	200	CF5	78.45		
		CF6	80.13		
	CF = (H/C)	CF7	69.168	%RSD = (Std Dev/Avg CF)*100	
		CF8	79.815		

**CONTINUING CALIBRATION A**

Date	7/8/2009		
Time	1750		
Instrument	A		
Compound	HMX		
Reported Conc	212	Rptd %D	6
Calculated Conc	211.653	Calcd %D	-5.83
Response	25610	ICAL CF	200
CF	121	CCV CF	211.7
	CF = (H/C)	%Difference = ((ICAL - CCV)/ICAL)*100	

**INITIAL CALIBRATION - C**

Date	6/9/2009	Rptd Avg CF	42.55812	Rptd %RSD	4.852
Instrument	C	Calcd Avg CF	42.558	Calcd %RSD	4.852
Compound	RDX				
Reported CF	42.585	CF1	45.40	STD DEV	2
Calculated CF	42.78	CF2	43.70		
		CF3	42.75		
Response	8556	CF4	43.30		
Conc	200	CF5	42.16		
		CF6	42.59		
	CF = (H/C)	CF7	38.118	%RSD = (Std Dev/Avg CF)*100	
		CF8	42.452		

**CONTINUING CALIBRATION C**

Date	7/10/2009		
Time	1932		
Instrument	C		
Compound	HMX		
Reported Conc	82.89	Rptd %D	17
Calculated Conc	82.894	Calcd %D	17.11
Response	3416	ICAL CF	100
CF	41.20913	CCV CF	82.9
	CF = (H/C)	%Difference = ((ICAL - CCV)/ICAL)*100	

**FIELD DUPLICATES**

COMPOUND	SAMPLE	QUALIFIER	DUPLICATE	QUALIFIER	RPD
PCA					#DIV/0!

NOTES      Samples are not qualified on this basis.

**COMPOUND IDENTIFICATION AND QUANTITATION**

Has a F10 been completed for every sample containing positive results?	Yes
Was RT data presented on the form?	NA
Are RTs within the established windows?	Yes
Any transcription or calculation errors?	No
Any false positives, negative peaks, shouldering, etc.?	No
Was GCMS confirmation needed for results > 10 µg/ml?	NA
Were percent differences or relative percent differences calculated?	NA
Are percent differences/RPDs greater than 25%?	NA
Are there any transcription errors?	No
Are Form 1s present for all field and quality control samples?	No
Are chromatograms and quan reports present for all samples?	Yes
Are RIs adjusted to reflect sample dilutions, percent solids, etc.?	Yes
For soils, any percent solids <50%?	No
For soils, any percent solids <10%?	No

23L4MW01AW RDX 0.13 µg/L

	A	C
response	524	269
cf	77.89163	42.55812
final vol ml	20	20
initial L	1.02157	1.02157
df	1	1
calculated	0.1317	0.1237

**PERCENT DIFFERENCE BETWEEN COLUMNS**

SAMPLE	COMPOUND	A	C	RPD	QUALIFIER
23L4MW01AW	RDX	0.1317	0.1237	6.3	none
23L4MW460W	HMX	0.028	0.0422	-40.5	none
	RDX	4.125	3.799	8.2	none
23L4MW03BW	HMX	0.0278	not confirmed	#VALUE!	none
	RDX	4.089	3.71	9.7	none
23L4MW03AW	HMX	0.3798	0.3908	-2.9	none
	RDX	9.369	8.736	7.0	none

LABORATORY CONTROL SAMPLE EVALUATION REPORT

HPLC

Client Lot #...: G9G010225      Work Order #...: LF1EN1AC      Matrix.....: WATER  
 LCS Lot-Sample#: G9G020000-251  
 Prep Date.....: 07/02/09      Analysis Date...: 07/08/09  
 Prep Batch #...: 9183251  
 Dilution Factor: 1

<u>PARAMETER</u>	<u>PERCENT RECOVERY</u>	<u>RECOVERY LIMITS</u>	<u>METHOD</u>
Nitroglycerin	98	(85 - 115)	SW846 8330
PETN	104	(84 - 117)	SW846 8330
Picric Acid	66	(21 - 118)	SW846 8330
2-Amino-4,6-dinitrotoluene	97	(77 - 123)	SW846 8330
4-Amino-2,6-dinitrotoluene	91	(68 - 113)	SW846 8330
1,3-Dinitrobenzene	98	(72 - 123)	SW846 8330
2,4-Dinitrotoluene	94	(70 - 119)	SW846 8330
2,6-Dinitrotoluene	95	(71 - 119)	SW846 8330
HMX	91	(67 - 115)	SW846 8330
Nitrobenzene	94	(69 - 119)	SW846 8330
2-Nitrotoluene	87	(64 - 120)	SW846 8330
3-Nitrotoluene	89	(67 - 114)	SW846 8330
4-Nitrotoluene	89	(67 - 115)	SW846 8330
RDX	97	(68 - 122)	SW846 8330
Tetryl	77	(66 - 105)	SW846 8330
1,3,5-Trinitrobenzene	106	(74 - 120)	SW846 8330
2,4,6-Trinitrotoluene	88	(69 - 111)	SW846 8330

<u>SURROGATE</u>	<u>PERCENT RECOVERY</u>	<u>RECOVERY LIMITS</u>
3,4-Dinitrotoluene	93	(79 - 111)

**NOTE (S) :**

Calculations are performed before rounding to avoid round-off errors in calculated results.

Bold print denotes control parameters

LABORATORY CONTROL SAMPLE DATA REPORT

HPLC

Client Lot #...: G9G010225      Work Order #...: LF1EN1AC      Matrix.....: WATER  
 LCS Lot-Sample#: G9G020000-251  
 Prep Date.....: 07/02/09      Analysis Date...: 07/08/09  
 Prep Batch #...: 9183251  
 Dilution Factor: 1

<u>PARAMETER</u>	<u>SPIKE</u> <u>AMOUNT</u>	<u>MEASURED</u> <u>AMOUNT</u>	<u>UNITS</u>	<u>PERCENT</u> <u>RECOVERY</u>	<u>METHOD</u>
Nitroglycerin	5.00	4.92	ug/L	98	SW846 8330
PTN	5.00	5.18	ug/L	104	SW846 8330
Picric Acid	5.00	3.28	ug/L	66	SW846 8330
2-Amino-4,6-dinitrotoluene	1.00	0.974	ug/L	97	SW846 8330
4-Amino-2,6-dinitrotoluene	1.00	0.909	ug/L	91	SW846 8330
1,3-Dinitrobenzene	1.00	0.976	ug/L	98	SW846 8330
2,4-Dinitrotoluene	1.00	0.936	ug/L	94	SW846 8330
2,6-Dinitrotoluene	1.00	0.946	ug/L	95	SW846 8330
HMX	1.00	0.914	ug/L	91	SW846 8330
Nitrobenzene	1.00	0.943	ug/L	94	SW846 8330
2-Nitrotoluene	1.00	0.869	ug/L	87	SW846 8330
3-Nitrotoluene	1.00	0.888	ug/L	89	SW846 8330
4-Nitrotoluene	1.00	0.892	ug/L	89	SW846 8330
RDX	1.00	0.970	ug/L	97	SW846 8330
Tetryl	1.00	0.772	ug/L	77	SW846 8330
1,3,5-Trinitrobenzene	1.00	1.06	ug/L	106	SW846 8330
2,4,6-Trinitrotoluene	1.00	0.880	ug/L	88	SW846 8330

<u>SURROGATE</u>	<u>PERCENT</u> <u>RECOVERY</u>	<u>RECOVERY</u> <u>LIMITS</u>
3,4-Dinitrotoluene	93	(79 - 111)

**NOTE(S):**

Calculations are performed before rounding to avoid round-off errors in calculated results.

Bold print denotes control parameters

METHOD BLANK REPORT

HPLC

Client Lot #...: G9G010225      Work Order #...: LF1EN1AA      Matrix.....: WATER  
 MB Lot-Sample #: G9G020000-251  
 Analysis Date...: 07/08/09      Prep Date.....: 07/02/09  
 Dilution Factor: 1              Prep Batch #...: 9183251

PARAMETER	RESULT	REPORTING		METHOD
		LIMIT	UNITS	
Nitroglycerin	ND	0.65	ug/L	SW846 8330
PETN	ND	0.65	ug/L	SW846 8330
Picric Acid	ND	1.0	ug/L	SW846 8330
2-Amino-4,6-dinitrotoluene	ND	0.20	ug/L	SW846 8330
4-Amino-2,6-dinitrotoluene	ND	0.10	ug/L	SW846 8330
1,3-Dinitrobenzene	ND	0.10	ug/L	SW846 8330
2,4-Dinitrotoluene	ND	0.10	ug/L	SW846 8330
2,6-Dinitrotoluene	ND	0.10	ug/L	SW846 8330
HMX	ND	0.10	ug/L	SW846 8330
Nitrobenzene	ND	0.10	ug/L	SW846 8330
2-Nitrotoluene	ND	0.40	ug/L	SW846 8330
3-Nitrotoluene	ND	0.40	ug/L	SW846 8330
4-Nitrotoluene	ND	0.50	ug/L	SW846 8330
RDX	ND	0.10	ug/L	SW846 8330
Tetryl	ND	0.10	ug/L	SW846 8330
1,3,5-Trinitrobenzene	ND	0.10	ug/L	SW846 8330
2,4,6-Trinitrotoluene	ND	0.10	ug/L	SW846 8330

SURROGATE	PERCENT	RECOVERY
	RECOVERY	LIMITS
3,4-Dinitrotoluene	94	(79 - 111)

NOTE(S):

Calculations are performed before rounding to avoid round-off errors in calculated results



A

Chromatography Summary

Injection Date: 7/8/2009 18:41 Operator: fhk  
 Data File: LC10 I07082009 BVA-000004.D Vial Num: 11  
 Instrument ID: LC10

Method 8330 Target Analyte Results

Sample: **LF1EN1AA 9183251 G9G020000-251**  
**B 1X**

Method File: LC10 I07082009.BV8330AB.M  
 Start Cal Date: 6/10/2009 15:33 End Cal Date: 6/11/2009 13:47

Dilution Factor	Extract Volume	Sample Volume	Sample Weight
1X	20 mL	1000 mL	0 g

Matrix: WATER SubList: WATER.sub SpikeList:  
 Samp. Info: LF1EN1AA 9183251 G9G020000-251 B 1X,0,  
 Misc. Info: ;,1000,,20;1;WATER sub,,0.1;LF1EN1AA

Signal 1 UV 250-265					Signal 2 UV 358-205					MDL	RL	Flag	
Compound Name	RT	Diff	Response	Conc (ug/L)	Flag	RT	Diff	Response	Conc (ug/L)	Flag	MDL	RL	Flag
3,4-Dinitrotoluene	18.23	0.041	5555	2.3550<		18.23	0.044	10715	2.4790		0.0000	0.00	
HMX											0.0270	0.10	
RDX											0.0650	0.10	
Picric ACID											0.1250	1.00	
1,3,5-Trinitrobenzene	10.37	0.184	1541	0.2088<							0.0310	0.10	45
1,3-Dinitrobenzene											0.0500	0.10	
TETRYL											0.0500	0.10	
Nitrobenzene											0.0500	0.10	
2,4,6-Trinitrotoluene											0.0240	0.10	
4-AM-2,6-DNT											0.0220	0.10	
2-AM-4,6-DNT											0.1000	0.20	
2,6-Dinitrotoluene											0.0500	0.10	
2,4-Dinitrotoluene											0.0500	0.10	
2-Nitrotoluene											0.0720	0.50	
4-Nitrotoluene											0.0720	0.50	
3-Nitrotoluene											0.0620	0.50	
Nitroglycerin											0.3300	0.65	
PETN											0.3000	0.65	
3,5-Dinitroaniline											0.0250	1.00	

*Not confirmed on W 10*  
*gmk 7/11/09*

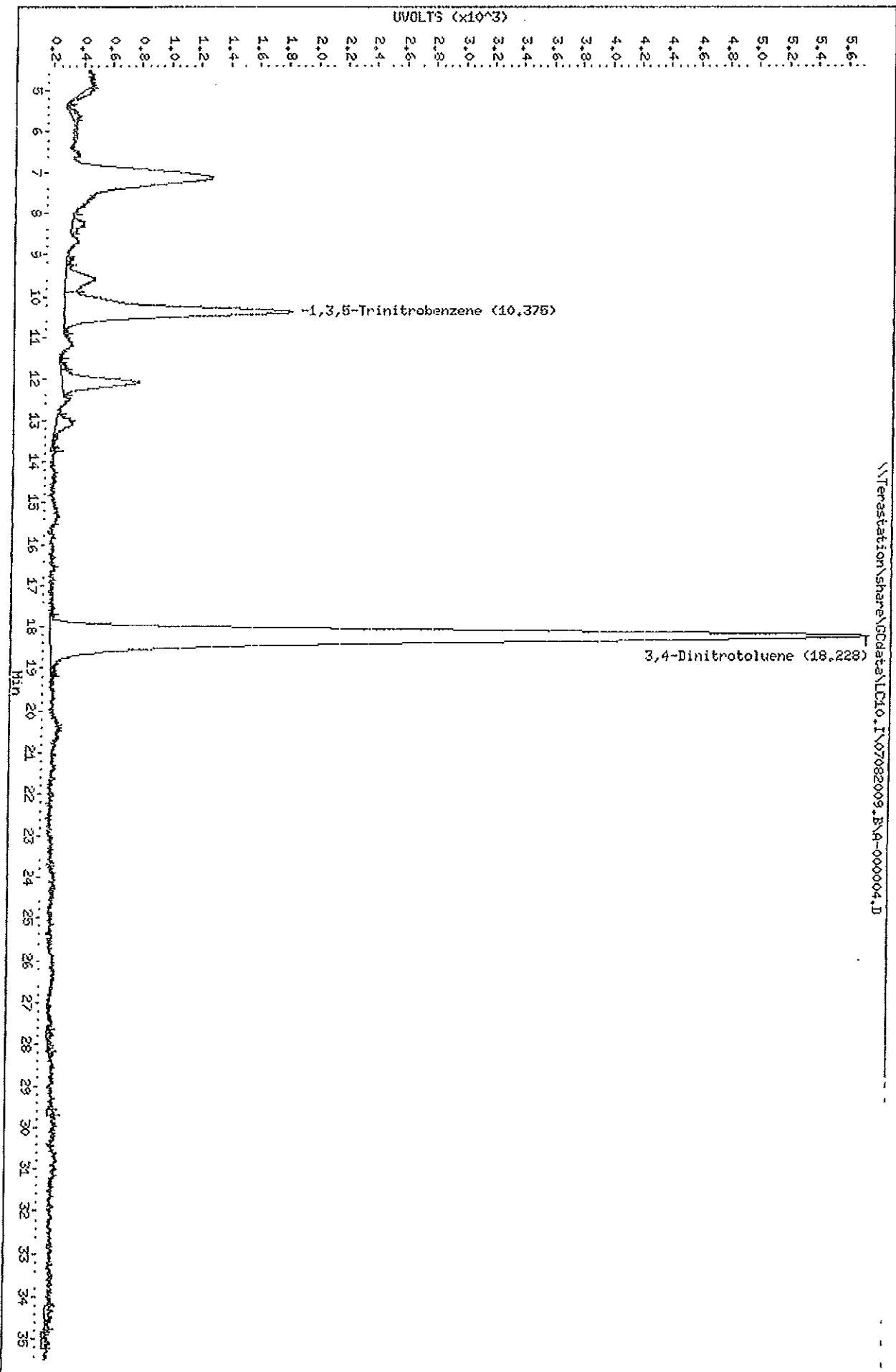
Surrogates:	Spiked	Recovered	%Rec	Spiked	Recovered	%Rec	Limits
3,4-Dinitrotoluene	2.5000	2.3550	94	2.5000	2.4790	99	(79-111)

Notes M = Manually Integrated Signals Differ by More Than 40%  
 D = Operator Disabled Result Signals Differ by More Than 50%  
 O = Over Calibration Range

Data File: \\Terastation\share\GCdata\LC10.I\07082009.B\A-000004.D  
Date : 08-JUL-2009 18:44

Client ID:  
Sample Info: LEIEN10A 9183251 09C020000-251 B 1X107  
Volume Injected (uL): 500.0  
Column phase: SYNERGI HYDRORP C18

Instrument: LC10.1  
Operator: PHK  
Column diameter: 4.60



Chromatography Summary

Injection Date: 7/10/2009 20:34 Operator: ftk  
 DataFile: LC9 107102009A BAC-000004 D Vial Num: 11  
 Instrument ID: LC9

Method 8330 Target Analyte Results

Method File: LC9 107102009A B8330CNAB M  
 Start Cal Date: 6/9/2009 20:53 End Cal Date: 6/10/2009 4:30

Sample: **LF1EN1AA 9183251 G9G020000-251**  
**B 1X**

Dilution Factor	Extract Volume	Sample Volume	Sample Weight
1X	20 mL	1000 mL	0 g

Matrix: WATER SubList: WATER sub SpikeList:  
 Samp. Info: LF1EN1AA 9183251 G9G020000-251 B 1X,0;  
 Misc. Info: ::1000;,20:1;WATER sub;,0,1;LF1EN1AA

Signal 1 UV 250-265						Signal 2 UV 358-205							
Compound Name	RT	Diff	Response	Conc (ug/L)	Flag	RT	Diff	Response	Conc (ug/L)	Flag	MDL	RL	Flag
3,4-Dinitrotoluene	30.78	0.093	5055	2.5500<							0.0000	0.00	45
HMX											0.0270	0.10	
RDX											0.0650	0.10	
Picric ACID											0.1250	1.00	
1,3,5-Trinitrobenzene											0.0310	0.10	
1,3-Dinitrobenzene											0.0500	0.10	
TETRYL											0.0500	0.10	
Nitrobenzene											0.0500	0.10	
2,4,6-Trinitrotoluene											0.0240	0.10	
4-AM-2,6-DNT											0.0220	0.10	
2-AM-4,6-DNT											0.1000	0.20	
2,6-Dinitrotoluene											0.0500	0.10	
2,4-Dinitrotoluene											0.0500	0.10	
2-Nitrotoluene											0.0720	0.50	
4-Nitrotoluene											0.0720	0.50	
3-Nitrotoluene											0.0620	0.50	
Nitroglycerin											0.3300	0.65	
PETN											0.3000	0.65	
3,5-Dinitroaniline											0.0250	1.00	

*Not confirmed MD  
 7/11/09*

Surrogates:	Spiked	Recovered	%Rec	Spiked	Recovered	%Rec	Limits
3,4-Dinitrotoluene	2.5000	2.5500	102	2.5000	0		(48-143)

Notes: M = Manually Integrated Signals Differ by More Than 40%  
 D = Operator Disabled Result Signals Differ by More Than 50%  
 O = Over Calibration Range

Data File: \\Terastation\share\SCdata\LC9.I\07102009A.B\C-000004.d  
Date: 10-JUL-2009 20:34

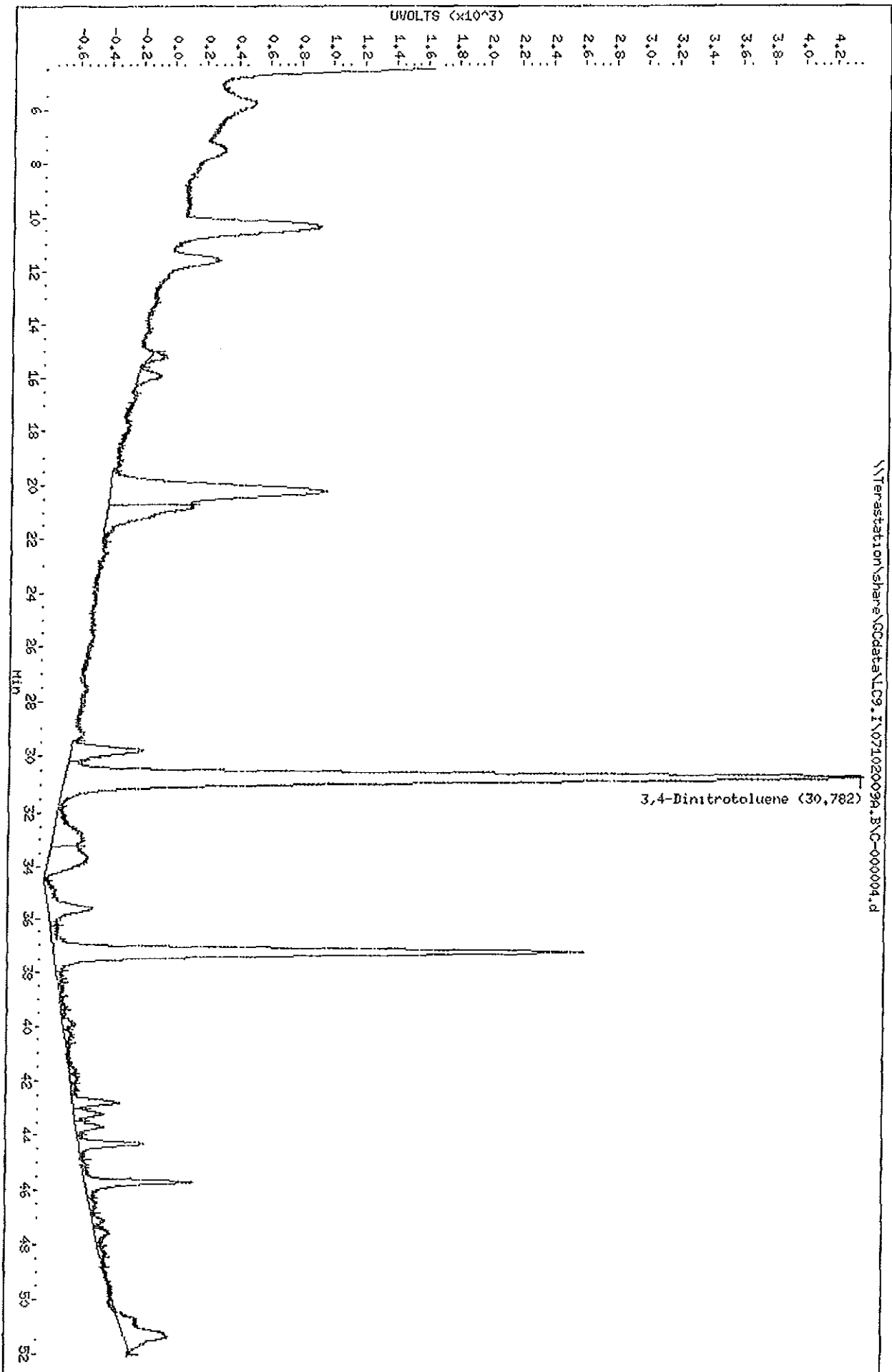
Client ID:

Sample Info: LFIEMHRA 9183251 696020000-251 B 1X10:  
Volume Injected (uL): 500.0  
Column phase: Agilent ZorbaxC18

Instrument: LC9.i

Operator: fmk  
Column diameter: 4.60

\\Terastation\share\SCdata\LC9.I\07102009A.B\C-000004.d



# QC DATA ASSOCIATION SUMMARY

G9G010225

Sample Preparation and Analysis Control Numbers

<u>SAMPLE#</u>	<u>MATRIX</u>	<u>ANALYTICAL METHOD</u>	<u>LEACH BATCH #</u>	<u>PREP BATCH #</u>	<u>MS RUN#</u>
001	WATER	SW846 8330		9183251	
002	WATER	SW846 8330		9183251	
003	WATER	SW846 8330		9183251	
004	WATER	SW846 8330		9183251	
005	WATER	SW846 8330		9183251	

TestAmerica West Sacramento  
ESC-Extraction Master Sheet

DE

Holding Time Due: 7-2-09 / 7-6-09  
BATCH #: 9183251

Project Due: 7-14-09  
Initiated By: MD  
Extn Comp'd By: TP

Date: 7-2-09  
Date: 7/8/09

Test #: 8330-L picnic & vid

QC Code	Lab ID	Sample Size	Initial Mass	Final Volume	Final Mass	Difference in Mass
B	MB	SW	Wt sheet			
C	LCS					
	696010225-01	7-2-09				
	696010225-02	MD				
	696010225-03					
	696010225-04					
	696010225-05					
	7/8/09 TP					

SOP No.: WLLI-0009

EXTRACTION COMMENTS:

Standard Information

QC Codes	Volume	STD ID	Exp. Date	STD Name/Conc	ppm/dpl
P11	50ul	096C5V0172	2-10-09	3240NT 50ul/ml	0.50
C	20ul	096C6V0037	7-17-09	8330-DNA 50ul/ml	0.50
BLIND	100ul	096C5V0149	11-4-09	NO PSTV 50ul/ml	1.00
BLANK	100ul	096C5V0039	7-17-09	PICNIC BLVD 50ul/ml	5.0

Spiked By / Date: MD - 7-2-09 Witnessed By / Date: TP 7/02/09

QC Codes: B = MB, C = LCS, L = LCSD, S = MS, D = SD, FB = Filter Blank

RCC058

TestAmerica Laboratories, Inc.  
EXTRACTION BENCH WORKSHEET

Run Date: 7/08/09  
Time: 10:59:23

LEV 1	LEV 2	Blank	LEV 1	LEV 2	Weights/Volumes
Y	Y	Check	Y	Y	Spike & Surrogate Worksheet
Y	Y	MS/MSD	Y	Y	Vial contains correct volume
Y	Y		Y	Y	Labels, Greenbars, worksheets
Y	Y		Y	Y	computer batch: correct & all match
Y	Y		Y	Y	Anomalies to Extraction Method

Expanded Deliverable  
COC Completed  
Bench Sheet Copied  
Package Submitted to Analytical Group  
Bench Sheet Copied per COC

\*\*\*\*\*  
\* QC BATCH: 9183251 \*  
\* PREP DATE: 7/02/09 9:30  
\* COMP DATE: 7/08/09 10:00  
\*\*\*\*\*

Extractionist: 000915 Horacio J. Arauz

Concentrationist: 002448 Tuan Q. Phan

Reviewer/Date: ARAUZH / 7/02/09

Nitroaromatics & Nitramines: Explosives (8330)  
SOLID PHASE EXTRACTION (NOMINAL)

EXTR EXPR	ANL DUE	LOT# WORK ORDER	TEST FLGS	EXT MTH	MATRIX	INIT/FIN WT/VOL	PH'S INIT ADJ1	ADJ2	EXTRACTION VOL	SOLVENTS VOL EXCHANGE	SPIKE STANDARD/SURROGATE ID
7/05/09	7/14/09	G9G010225-001 LFXKX-1-AC	R B7 A0	A0	WATER	1021.57mL 20.00mL	NA	NA	HOAC/ACN	4.5 HOAC/ACN	5.0 50UL-09GCSV0172
7/06/09	7/14/09	G9G010225-002 LFXK6-1-AC	R B7 A0	A0	WATER	1022.4mL 20.00mL	NA	NA	HOAC/ACN	4.5 HOAC/ACN	5.0 50UL-09GCSV0172
7/03/09	7/14/09	G9G010225-003 LFX11-1-AC	R B7 A0	A0	WATER	1024.15mL 20.00mL	NA	NA	HOAC/ACN	4.5 HOAC/ACN	5.0 50UL-09GCSV0172
7/03/09	7/14/09	G9G010225-004 LFX12-1-AC	R B7 A0	A0	WATER	1019.2mL 20.00mL	NA	NA	HOAC/ACN	4.5 HOAC/ACN	5.0 50UL-09GCSV0172
7/03/09	7/14/09	G9G010225-005 LFX15-1-AC	R B7 A0	A0	WATER	1021.92mL 20.00mL	NA	NA	HOAC/ACN	4.5 HOAC/ACN	5.0 50UL-09GCSV0172
7/06/09	0/00/00	G9G020000-251 LFX1N-1-AAB	B7 A0	A0	WATER	1000mL 20.00mL	NA	NA	HOAC/ACN	4.5 HOAC/ACN	5.0 50UL-09GCSV0172
7/06/09	0/00/00	G9G020000-251 LFX1N-1-ACC	B7 A0	A0	WATER	1000mL 20.00mL	NA	NA	HOAC/ACN	4.5 HOAC/ACN	5.0 50UL-09GCSV0172

100UL-09GCSV0149; 100UL-09GCSV0039, SEE COLUMN WATER 003339096A;  
-45 FILTER MILLEPORE LOT RSPN29184; SODIUM CHLORIDE MALLINCKRODT LOT H07582  
-1% HOAC/ACN 2991-95F; 1% HOAC/H2O 2991-87B

R = RUSH C = CLP  
E = EPA 600 D = EXP.DEL)

NUMBER OF WORK ORDERS IN BATCH: 7

A

TestAmerica West Sacramento  
GC/LC INSTRUMENT LOG

Inst ID: LC10                      Batch ID: 06102009  
Method : Method 8330              Test : SOP SAC-LC-0009  
ICAL Date: See Calibration Report

Date	Time	Operator	Sample ID	File ID	Vol or Wt	Extract Vol	Diln	Comments
10-JUN-2009	14:42	fhk	Blank	A-000001.	0 g	0 mL	1	
10-JUN-2009	15:33	fhk	STD_1 09GCSV0048 5.0/0/0/0ng/mL	A-000002.	0 g	0 mL	1	
10-JUN-2009	16:23	fhk	STD_2 09GCSV0049 10/20/10/10ng	A-000003.	0 g	0 mL	1	
10-JUN-2009	17:14	fhk	STD_3 09GCSV0050 20/50/20/20ng	A-000004.	0 g	0 mL	1	Bad Std
10-JUN-2009	18:05	fhk	STD_4 09GCSV0051 50/100/50/50ng	A-000005.	0 g	0 mL	1	
10-JUN-2009	18:56	fhk	STD_5 09GCSV0053 100/200/100/100ng	A-000006.	0 g	0 mL	1	
10-JUN-2009	19:46	fhk	STD_6 09GCSV0054 200/500/200/200ng	A-000007.	0 g	0 mL	1	
10-JUN-2009	20:37	fhk	STD_7 09GCSV0055 500/1000/500/1000ng	A-000008.	0 g	0 mL	1	
10-JUN-2009	21:28	fhk	STD_8 09GCSV0056 1000/2000/1000/1000ng	A-000009.	0 g	0 mL	1	
10-JUN-2009	22:18	fhk	Blank	A-000010.	0 g	0 mL	1	
10-JUN-2009	23:09	fhk	ICV_6 08GCSV0397 100/200/100/100ng	A-000011.	0 g	0 mL	1	
11-JUN-2009	00:00	fhk	STD_5 09GCSV0053 100/200/100/100ng	A-000012.	0 g	0 mL	1	
11-JUN-2009	00:50	fhk	Surrogate 100ng/mL	A-000013.	0 g	0 mL	1	
11-JUN-2009	12:56	fhk	Primer	A-000014.	0 g	0 mL	1	
11-JUN-2009	13:47	fhk	STD_3 09GCSV0050 20/50/20/20ng	A-000015.	0 g	0 mL	1	best select / new Std

John 6/11/09



TestAmerica West Sacramento

INITIAL CALIBRATION DATA

Start Cal Date : 10-JUN-2009 15:33  
 End Cal Date : 11-JUN-2009 13:47  
 Quant Method : ESTD  
 Origin : Disabled  
 Target Version : 4.14  
 Integrator : Falcon  
 Method file : \\Terastation\share\GCdata\LC10.I\06102009.B\8330AB.M  
 Last Edit : 11-Jun-2009 15:06 kenneyf  
 Curve Type : Average

Calibration File Names:

- Level 1: \\Terastation\share\GCdata\GCdata\LC10.I\06102009.B\A-000002.d
- Level 2: \\Terastation\share\GCdata\GCdata\LC10.I\06102009.B\A-000003.d
- Level 3: \\Terastation\share\GCdata\GCdata\LC10.I\06102009.B\A-000015.d
- Level 4: \\Terastation\share\GCdata\GCdata\LC10.I\06102009.B\A-000005.d
- Level 5: \\Terastation\share\GCdata\GCdata\LC10.I\06102009.B\A-000006.d
- Level 6: \\Terastation\share\GCdata\GCdata\LC10.I\06102009.B\A-000007.d
- Level 7: \\Terastation\share\GCdata\GCdata\LC10.I\06102009.B\A-000008.d
- Level 8: \\Terastation\share\GCdata\GCdata\LC10.I\06102009.B\A-000009.d

Compound	5.000	10.000	20.000	50.000	100.000	200.000	RRF	% RSD
	Level 1	Level 2	Level 3	Level 4	Level 5	Level 6		
	500.000	1000.000						
	Level 7	Level 8						
2 HMX	119 ✓ 116 ✓	120 ✓ 123 ✓	115 ✓	125 ✓	123 ✓	124 ✓	121	3.066
3 RDX	81.80000 69.16600	75.90000 ✓ 79.81500 ✓	77.15000 ✓	80.77000 ✓	78.45000 ✓	80.13000 ✓	77.89163	5.156
4 EGDN	++++ +++	++++ ++++	++++	++++	++++	++++	++++	++++
5 Picric ACID	++++ 74.63400	++++ 74.87800	91.72000 ✓	85.68000 ✓	84.85500 ✓	84.88200 ✓	82.77483	8.120
6 1,3,5-Trinitrobenzene	161 136	146 147	145	150	147	148	148	4.610
7 1,3-Dinitrobenzene	145 128	142 141	138	142	139	142	140	3.715
8 3,5-Dinitroaniline	97.20000 82.20800	90.80000 89.25000	89.35000	91.80000	88.83000	90.35000	89.97350	4.581
9 TETRYL	97.60000 92.43000	95.10000 ✓ 94.40900 ✓	77.40000 ✓	96.02000 ✓	94.31000 ✓	95.06000 ✓	92.79113	6.888

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TestAmerica West Sacramento

INITIAL CALIBRATION DATA

Start Cal Date : 10-JUN-2009 15:33  
 End Cal Date : 11-JUN-2009 13:47  
 Quant Method : ESTD  
 Origin : Disabled  
 Target Version : 4.14  
 Integrator : Falcon  
 Method file : \\Terastation\share\GCdata\LC10.I\06102009.B\8330AB.M  
 Last Edit : 11-Jun-2009 15:06 kenneyf  
 Curve Type : Average

Compound	5.000	10.000	20.000	50.000	100.000	200.000	RRF	% RSD
	Level 1	Level 2	Level 3	Level 4	Level 5	Level 6		
	500.000	1000.000						
	Level 7	Level 8						
10 Nitrobenzene	63.80000	62.30000	56.55000	61.42000	61.06000	62.26000		
	56.02600	61.95700					60.67412	4.659
11 Nitroglycerin	+++++	+++++	+++++	+++++	+++++	+++++	+++++	+++++
12 2,4,6-Trinitrotoluene	94.80000	90.20000	81.90000	88.78000	87.04000	87.55500		
	84.22200	87.17800					87.70937	4.401
13 4-AM-2,6-DNT	67.00000	63.30000	60.85000	63.26000	61.61000	62.67000		
	58.65400	61.14600					62.31125	3.901
14 2-AM-4,6-DNT	74.80000	75.80000	73.10000	75.54000	73.53000	74.39000		
	69.18400	72.13500					73.55987	2.932
15 2,6-Dinitrotoluene	57.60000	53.30000	51.90000	53.68000	53.03000	53.21500		
	50.43000	52.74900					53.23800	3.835
16 2,4-Dinitrotoluene	93.00000	86.50000	82.60000	86.20000	84.72000	85.49000		
	81.43200	85.05700					85.62488	4.031
17 2-Nitrotoluene	40.80000	37.40000	33.20000	35.64000	35.21000	35.77500		
	33.97400	35.45000					35.93113	6.486
18 4-Nitrotoluene	48.00000	46.40000	40.75000	43.56000	43.22000	43.56500		
	41.67600	43.21300					43.79800	5.390
19 3-Nitrotoluene	47.00000	44.60000	39.85000	43.24000	42.93000	43.08000		
	41.51600	42.82400					43.13000	4.852
20 PETN	+++++	+++++	+++++	+++++	+++++	+++++	+++++	+++++

TestAmerica West Sacramento

INITIAL CALIBRATION DATA

Start Cal Date : 10-JUN-2009 15:33  
 End Cal Date : 11-JUN-2009 13:47  
 Quant Method : ESTD  
 Origin : Disabled  
 Target Version : 4.14  
 Integrator : Falcon  
 Method file : \\Terastation\share\GCdata\LC10.I\06102009.B\8330AB.M  
 Last Edit : 11-Jun-2009 15:06 kenneyf  
 Curve Type : Average

Compound	5.000	10.000	20.000	50.000	100.000	200.000	RRP	% RSD
	Level 1	Level 2	Level 3	Level 4	Level 5	Level 6		
	500.000	1000.000						
	Level 7	Level 8						
\$ 1 3,4-Dinitrotoluene	+++++	50.10000	46.40000	46.48000	45.87000	46.85500		
	46.08000	48.38600					47.16729	3.251

TestAmerica West Sacramento

INITIAL CALIBRATION DATA

Start Cal Date : 10-JUN-2009 16:23  
 End Cal Date : 11-JUN-2009 13:47  
 Quant Method : ESTD  
 Origin : Disabled  
 Target Version : 4.14  
 Integrator : Falcon  
 Method file : \\Terastation\share\GCdata\LC10.I\06102009.B\8330AB.M\81  
 Last Edit : 11-Jun-2009 15:14 kenneyf  
 Curve Type : Average

Calibration File Names:

Level 2: \\Terastation\share\GCdata\GCdata\LC10.I\06102009.B\A-000003.d\A.  
 Level 3: \\Terastation\share\GCdata\GCdata\LC10.I\06102009.B\A-000015.d\A.  
 Level 4: \\Terastation\share\GCdata\GCdata\LC10.I\06102009.B\A-000005.d\A.  
 Level 5: \\Terastation\share\GCdata\GCdata\LC10.I\06102009.B\A-000006.d\A.  
 Level 6: \\Terastation\share\GCdata\GCdata\LC10.I\06102009.B\A-000007.d\A.  
 Level 7: \\Terastation\share\GCdata\GCdata\LC10.I\06102009.B\A-000008.d\A.  
 Level 8: \\Terastation\share\GCdata\GCdata\LC10.I\06102009.B\A-000009.d\A.

Compound	10.000	20.000	50.000	100.000	200.000	500.000	RRF	% RSD
	Level 2	Level 3	Level 4	Level 5	Level 6	Level 7		
	1000.000							
	Level 8							
2 HMX	++++	++++	++++	++++	++++	++++	++++	++++
3 RDX	++++	++++	++++	++++	++++	++++	++++	++++
4 EGDN	++++	++++	++++	++++	++++	++++	++++	++++
5 Picric ACID	++++	134	126	125	125	110	122	8.080
6 1,3,5-Trinitrobenzene	++++	++++	++++	++++	++++	++++	++++	++++
7 1,3-Dinitrobenzene	++++	++++	++++	++++	++++	++++	++++	++++
8 3,5-Dinitroaniline	++++	++++	++++	++++	++++	++++	++++	++++
9 TETRYL	++++	++++	++++	++++	++++	++++	++++	++++

TestAmerica West Sacramento

INITIAL CALIBRATION DATA

Start Cal Date : 10-JUN-2009 16:23  
 End Cal Date : 11-JUN-2009 13:47  
 Quant Method : ESTD  
 Origin : Disabled  
 Target Version : 4.14  
 Integrator : Falcon  
 Method file : \\Terastation\share\GCdata\LC10.I\06102009.B\8330AB.M\83  
 Last Edit : 11-Jun-2009 15:14 kenneyf  
 Curve Type : Average

Compound	10.000 Level 2	20.000 Level 3	50.000 Level 4	100.000 Level 5	200.000 Level 6	500.000 Level 7	RRF	% RSD
	1000.000 Level 8							
10 Nitrobenzene	+++++	+++++	+++++	+++++	+++++	+++++	+++++	+++++
11 Nitroglycerin	+++++	61.25000	59.82000	60.58000	61.86500	59.75400	60.83467	1.527
12 2,4,6-Trinitrotoluene	+++++	+++++	+++++	+++++	+++++	+++++	+++++	+++++
13 4-AM-2,6-DNT	+++++	+++++	+++++	+++++	+++++	+++++	+++++	+++++
14 2-AM-4,6-DNT	+++++	+++++	+++++	+++++	+++++	+++++	+++++	+++++
15 2,6-Dinitrotoluene	+++++	+++++	+++++	+++++	+++++	+++++	+++++	+++++
16 2,4-Dinitrotoluene	+++++	+++++	+++++	+++++	+++++	+++++	+++++	+++++
17 2-Nitrotoluene	+++++	+++++	+++++	+++++	+++++	+++++	+++++	+++++
18 4-Nitrotoluene	+++++	+++++	+++++	+++++	+++++	+++++	+++++	+++++
19 3-Nitrotoluene	+++++	+++++	+++++	+++++	+++++	+++++	+++++	+++++
20 PETN	+++++	32.65000	30.36000	31.91000	32.25000	32.38000	32.05033	2.748

TestAmerica West Sacramento

INITIAL CALIBRATION DATA

Start Cal Date : 10-JUN-2009 16:23  
 End Cal Date : 11-JUN-2009 13:47  
 Quant Method : ESTD  
 Origin : Disabled  
 Target Version : 4.14  
 Integrator : Falcon  
 Method file : \\Terastation\share\GCdata\LC10.I\06102009.B\8330AB.M\83  
 Last Edit : 11-Jun-2009 15:14 kenneyf  
 Curve Type : Average

Compound	10.000 Level 2	20.000 Level 3	50.000 Level 4	100.000 Level 5	200.000 Level 6	500.000 Level 7	RRF	% RSD
	1000.000 Level 8							
\$ 1,3,4-Dinitrotoluene	*++++ 93.14200	90.40000	88.76000	88.28000	90.51000	88.03667	89.85478	2.142

*\* Level 2 dropped due to poor integration  
 JMK  
 6/12/09*

Chromatography Summary

Injection Date: 6/10/2009 23:09 Operator: fhk  
 DataFile: LC10 N05102009 BVA-000011.D Vial Num: 20  
 Instrument ID: LC10

Method 8330 Target Analyte Results

Sample: **ICV\_6 08GCSV0397**  
**100/200/100/100ng/mL**

Method File: LC10 N05102009.BW8330AB.M  
 Start Cal Date: 6/10/2009 15:33 End Cal Date: 6/11/2009 13:47

Matrix: NONE SubList: CAL.sub SpikeList:  
 Samp. Info: ICV\_6 08GCSV0397 100/200/100/100ng/mL:2  
 Misc. Info: ;6; ; ;3;CAL.sub; ;0;1

Dilution Factor	Extract Volume	Sample Volume	Sample Weight
1X	0 mL	0 mL	0 g

Signal 1 UV 250-265								Signal 2 UV 358-205								
Compound Name	RT	Response	PPB	Spike Level	%D	Result	Flag	RT	Response	PPB	Spike Level	%D	Result	Flag	Limits(%)	Flag
<b>3,4-Dinitrotoluene</b>				<b>200</b>	<b>-100%</b>	<b>Fails</b>					<b>200</b>	<b>-100%</b>	<b>Fails</b>		(±15)	
HMX	5.29	26153	216.4000<	200	8%	Acceptable					200	-100%	Fails		(±15)	45
RDX	7.84	16084	205.5000<	200	3%	Acceptable					200	-100%	Fails		(±15)	45
Picric ACID	9.00	41754	504.4000 ✓	500	1%	Acceptable		9.00	61509	505.5000<	500	1%	Acceptable		(±15)	
1,3,5-Trinitrobenzene	10.20	29097	197.1000<	200	-1%	Acceptable					200	-100%	Fails		(±15)	45
1,3-Dinitrobenzene	13.25	27834	199.2000<	200	0%	Acceptable					200	-100%	Fails		(±15)	45
TETRYL	14.30	17776	191.6000<	200	-4%	Acceptable					200	-100%	Fails		(±15)	45
Nitrobenzene	15.16	12143	200.1000<	200	0%	Acceptable					200	-100%	Fails		(±15)	45
2,4,6-Trinitrotoluene	16.53	17352	197.8000<	200	-1%	Acceptable					200	-100%	Fails		(±15)	45
4-AM-2,6-DNT	17.51	11969	192.1000< ✓	200	-4%	Acceptable					200	-100%	Fails		(±15)	45
2-AM-4,6-DNT	18.57	13851	188.3000<	200	-6%	Acceptable					200	-100%	Fails		(±15)	45
2,6-Dinitrotoluene	20.28	10445	196.2000<	200	-2%	Acceptable					200	-100%	Fails		(±15)	45
2,4-Dinitrotoluene	21.00	16653	194.5000<	200	-3%	Acceptable					200	-100%	Fails		(±15)	45
2-Nitrotoluene	24.68	7059	196.7000<	200	-2%	Acceptable					200	-100%	Fails		(±15)	45
4-Nitrotoluene	26.50	8526	194.7000<	200	-3%	Acceptable					200	-100%	Fails		(±15)	45
3-Nitrotoluene	28.49	8451	195.9000< ✓	200	-2%	Acceptable					200	-100%	Fails		(±15)	45
Nitroglycerin				<b>200</b>	<b>-100%</b>	<b>Fails</b>		15.78	12548	206.3000< ✓	200	3%	Acceptable		(±15)	45
PETN				<b>200</b>	<b>-100%</b>	<b>Fails</b>		30.10	5830	181.9000< ✓	200	-9%	Acceptable		(±15)	45
3,5-Dinitroaniline	13.99	18460	205.2000<	200	3%	Acceptable					200	-100%	Fails		(±15)	45
EGDN				<b>200</b>	<b>-100%</b>	<b>Fails</b>					200	-100%	Fails		(±15)	

Notes: M = Manually Integrated Signals Differ by More Than 40%  
 D = Operator Disabled Result Signals Differ by More Than 50%  
 O = Over Calibration Range





Chromatography Summary

Injection Date: 7/8/2009 17:50 Operator: ftk  
 DataFile: I:\10\07082009\BVA-000003.D Vial Num: 2  
 Instrument ID: I.C10

Method 8330 Target Analyte Results

Sample: STD\_6 09GCSV0054 .4K/2/2/2/2

Method File: I:\10\07082009\BVA8330AB.M  
 Start Cal Date: 6/10/2009 15:33 End Cal Date: 6/11/2009 13:47

Matrix: NONE SubList: CAL sub SpikeList:  
 Samp. Info: STD\_6 09GCSV0054 .4K/2/2/2/2  
 Misc. Info: .6 . . . 3,CAL sub, .0,1

Dilution Factor	Extract Volume	Sample Volume	Sample Weight
IX	0 mL	0 mL	0 g

Signal 1 UV 250-265								Signal 2 UV 358-205								
Compound Name	RT	Response	PPB	Spike Level	%D	Result	Flag	RT	Response	PPB	Spike Level	%D	Result	Flag	Limits(%)	Flag
3,4-Dinitrotoluene	18.19	9699	205.6000<	200	3%	Acceptable		18.18	18512	214.1000	200	7%	Acceptable		(±15)	
HMX	5.28	25610	212.0000<	200	6%	Acceptable					200	-100%	Fails		(±15)	45
RDX	7.84	16375	210.2000<	200	5%	Acceptable					200	-100%	Fails		(±15)	45
Picric ACID	8.99	41964	507.0000	500	1%	Acceptable		8.99	61727	507.3000<	500	1%	Acceptable		(±15)	
1,3,5-Trinitrobenzene	10.19	30519	206.7000<	200	3%	Acceptable					200	-100%	Fails		(±15)	45
1,3-Dinitrobenzene	13.25	29220	209.2000<	200	5%	Acceptable					200	-100%	Fails		(±15)	45
TETRYL	14.32	19026	205.0000<	200	3%	Acceptable					200	-100%	Fails		(±15)	45
Nitrobenzene	15.17	12927	213.0000<	200	7%	Acceptable					200	-100%	Fails		(±15)	45
2,4,6-Trinitrotoluene	16.54	18133	206.7000<	200	3%	Acceptable					200	-100%	Fails		(±15)	45
4-AM-2,6-DNT	17.53	12950	207.8000<	200	4%	Acceptable					200	-100%	Fails		(±15)	45
2-AM-4,6-DNT	18.59	15167	206.2000<	200	3%	Acceptable					200	-100%	Fails		(±15)	45
2,6-Dinitrotoluene	20.32	10979	206.2000<	200	3%	Acceptable					200	-100%	Fails		(±15)	45
2,4-Dinitrotoluene	21.04	17596	205.5000<	200	3%	Acceptable					200	-100%	Fails		(±15)	45
2-Nitrotoluene	24.75	7372	205.2000<	200	3%	Acceptable					200	-100%	Fails		(±15)	45
4-Nitrotoluene	26.57	8973	204.9000<	200	2%	Acceptable					200	-100%	Fails		(±15)	45
3-Nitrotoluene	28.57	8900	206.4000<	200	3%	Acceptable					200	-100%	Fails		(±15)	45
Nitroglycerin				200	-100%	Fails		15.79	12780	210.1000<	200	5%	Acceptable		(±15)	45
PIETN				200	-100%	Fails		30.20	6780	211.5000<	200	6%	Acceptable		(±15)	45
3,5-Dinitroaniline	13.99	18618	206.9000<	200	3%	Acceptable					200	-100%	Fails		(±15)	45
EGDN				200	-100%	Fails					200	-100%	Fails		(±15)	

Notes: M = Manually Integrated Signals Differ by More Than 40%  
 D = Operator Disabled Result Signals Differ by More Than 50%  
 O = Over Calibration Range

Chromatography Summary

Injection Date: 7/9/2009 0:35 Operator: ftk  
 Data File: LC10 N07082009.BVA-000011.D Vial Num: 3  
 Instrument ID: LC10

Method 8330 Target Analyte Results

Sample: **STD\_5 09GCSV0053 .2K/1/1/1/1**

Method File: LC10 N07082009 B8330AB.M  
 Start Cal Date: 6/10/2009 15:33 End Cal Date: 6/11/2009 13:47

Matrix: NONE SubList: CAL.sub SpikeList:  
 Samp. Info: STD\_5 09GCSV0053 .2K/1/1/1/1;2  
 Misc. Info: ;5; ; ;3.CAL.sub: 0.1

Dilution Factor	Extract Volume	Sample Volume	Sample Weight
1X	0 mL	0 mL	0 g

Signal 1 UV 250-265							Signal 2 UV 358-205							
Compound Name	RT	Response	PPB	Spike Level	%D	Result Flag	RT	Response	PPB	Spike Level	%D	Result Flag	Limits(%)	Flag
3,4-Dinitrotoluene	18.20	4739	100 5000<	100	1%	Acceptable	18.21	9063	104 8000	100	5%	Acceptable	(±15)	
HMX	5.29	12630	104 5000<	100	5%	Acceptable				100	-100%	Fails	(±15)	45
RDX	7.85	8029	103 1000<	100	3%	Acceptable				100	-100%	Fails	(±15)	45
Picric ACID	9.05	16777	202 7000	200	1%	Acceptable	9.06	24679	202 8000<	200	1%	Acceptable	(±15)	
1,3,5-Trinitrobenzene	10.20	15097	102 3000<	100	2%	Acceptable				100	-100%	Fails	(±15)	45
1,3-Dinitrobenzene	13.26	14436	103 3000<	100	3%	Acceptable				100	-100%	Fails	(±15)	45
TETRYL	14.34	9407	101 4000<	100	1%	Acceptable				100	-100%	Fails	(±15)	45
Nitrobenzene	15.18	6290	103 7000<	100	4%	Acceptable				100	-100%	Fails	(±15)	45
2,4,6-Trinitrotoluene	16.56	8976	102 3000<	100	2%	Acceptable				100	-100%	Fails	(±15)	45
4-AM-2,6-DNT	17.55	6365	102 1000<	100	2%	Acceptable				100	-100%	Fails	(±15)	45
2-AM-4,6-DNT	18.62	7472	101 6000<	100	2%	Acceptable				100	-100%	Fails	(±15)	45
2,6-Dinitrotoluene	20.35	5434	102 1000<	100	2%	Acceptable				100	-100%	Fails	(±15)	45
2,4-Dinitrotoluene	21.06	8697	101 6000<	100	2%	Acceptable				100	-100%	Fails	(±15)	45
2-Nitrotoluene	24.79	3631	101 0000<	100	1%	Acceptable				100	-100%	Fails	(±15)	45
4-Nitrotoluene	26.60	4405	100 6000<	100	1%	Acceptable				100	-100%	Fails	(±15)	45
3-Nitrotoluene	28.62	4379	101 5000<	100	2%	Acceptable				100	-100%	Fails	(±15)	45
Nitroglycerin				100	-100%	Fails	15.81	6320	103 9000<	100	4%	Acceptable	(±15)	45
PETN				100	-100%	Fails	30.24	3380	105 4000<	100	5%	Acceptable	(±15)	45
3,5-Dinitroaniline	14.01	9130	101 5000<	100	2%	Acceptable				100	-100%	Fails	(±15)	45
EGDN				100	-100%	Fails				100	-100%	Fails	(±15)	

Notes: M = Manually Integrated Signals Differ by More Than 40%  
 D = Operator Disabled Result Signals Differ by More Than 50%  
 O = Over Calibration Range



TestAmerica West Sacramento

INITIAL CALIBRATION DATA

Start Cal Date : 09-JUN-2009 20:53  
 End Cal Date : 10-JUN-2009 04:20  
 Quant Method : ESTD  
 Origin : Disabled  
 Target Version : 4.14  
 Integrator : HP Genie  
 Method file : \\Terastation\share\GCdata\LC9.I\06092009.B\8330CNAB.M  
 Last Edit : 10-Jun-2009 10:24 kenneyf  
 Curve Type : Average

Calibration File Names:

- Level 1: \\Terastation\share\GCdata\GCdata\LC9.I\06092009.B\C-000004.d
- Level 2: \\Terastation\share\GCdata\GCdata\LC9.I\06092009.B\C-000005.d
- Level 3: \\Terastation\share\GCdata\GCdata\LC9.I\06092009.B\C-000006.d
- Level 4: \\Terastation\share\GCdata\GCdata\LC9.I\06092009.B\C-000007.d
- Level 5: \\Terastation\share\GCdata\GCdata\LC9.I\06092009.B\C-000008.d
- Level 6: \\Terastation\share\GCdata\GCdata\LC9.I\06092009.B\C-000009.d
- Level 7: \\Terastation\share\GCdata\GCdata\LC9.I\06092009.B\C-000010.d
- Level 8: \\Terastation\share\GCdata\GCdata\LC9.I\06092009.B\C-000011.d

Compound	5.000	10.000	20.000	50.000	100.000	200.000	RRF	% RSD
	Level 1	Level 2	Level 3	Level 4	Level 5	Level 6		
	500.000	1000.000						
	Level 7	Level 8						
2 HMX	49.60000 37.41800	43.70000 39.22000	40.45000	40.46000	39.51000	39.31500	41.20913	9.288
3 RDX	45.40000 38.11800	43.70000 42.45200	42.75000	43.30000	42.16000	42.58500	42.55812	4.652
4 EGDN	++++	++++	++++	++++	++++	++++	++++	++++
5 Picric ACID	++++	++++	++++	++++	++++	++++	++++	++++
6 1,3,5-Trinitrobenzene	66.80000 63.46600	67.20000 67.27500	67.60000	68.82000	67.02000	67.61500	66.97450	2.306
7 1,3-Dinitrobenzene	95.20000 79.37000	88.70000 88.38800	87.25000	89.40000	87.58000	88.56500	88.05662	4.893
8 3,5-Dinitroaniline	71.20000 60.97200	68.50000 65.89900	68.65000	68.54000	66.12000	66.97500	67.10700	4.478
9 TETRYL	93.20000 85.04400	86.50000 86.33600	85.70000	86.92000	86.05000	86.26500	87.00187	2.950

TestAmerica West Sacramento

INITIAL CALIBRATION DATA

Start Cal Date : 09-JUN-2009 20:53  
 End Cal Date : 10-JUN-2009 04:20  
 Quant Method : ESTD  
 Origin : Disabled  
 Target Version : 4.14  
 Integrator : HP Genie  
 Method file : \\Terastation\share\GCdata\LC9.I\06092009.B\8330CNAB.M  
 Last Edit : 10-Jun-2009 10:24 kenneyf  
 Curve Type : Average

Compound	5.000	10.000	20.000	50.000	100.000	200.000	RRF	% RSD
	Level 1	Level 2	Level 3	Level 4	Level 5	Level 6		
	500.000	1000.000						
	Level 7	Level 8						
10 Nitrobenzene	39.40000 35.35200	42.00000 40.30500	38.65000	40.64000	39.79000	40.66500	39.60025	5.005
11 Nitroglycerin	++++ ++++	++++ ++++	++++	++++	++++	++++	++++	++++
12 2,4,6-Trinitrotoluene	104 71.97000	91.00000 73.60800	81.00000	77.16000	74.32000	73.90500	80.89538	13.892
13 4-AM-2,6-DNT	68.80000 64.80800	69.80000 65.14700	68.05000	57.74000	65.37000	65.85500	66.94625	2.818
14 2-AM-4,6-DNT	80.20000 69.79400	76.80000 72.19800	75.45000	74.54000	72.25000	73.07500	74.28838	4.343
15 2,6-Dinitrotoluene	54.80000 45.69200	52.40000 49.41100	49.70000	49.88000	49.11000	49.44000	50.05412	5.280
16 2,4-Dinitrotoluene	83.00000 73.10400	82.40000 78.73900	79.60000	30.22000	78.86000	79.32000	79.40538	3.779
17 2-Nitrotoluene	24.50000 22.18600	25.30000 24.28200	23.52500	24.59000	24.32500	24.51250	24.12756	3.813
18 4-Nitrotoluene	++++ ++++	++++ ++++	++++	++++	++++	++++	++++	++++
19 3-Nitrotoluene	35.00000 28.82000	32.50000 30.96400	30.55000	31.42000	30.76000	31.17500	31.39862	5.672
20 PETN	++++ ++++	++++ ++++	++++	++++	++++	++++	++++	++++

TestAmerica West Sacramento

INITIAL CALIBRATION DATA

Start Cal Date : 09-JUN-2009 20:53  
 End Cal Date : 10-JUN-2009 04:20  
 Quant Method : ESTD  
 Origin : Disabled  
 Target Version : 4.14  
 Integrator : HP Genie  
 Method file : \\Terastation\share\GCdata\LC9.I\06092009.B\8330CNAB.M  
 Last Edit : 10-Jun-2009 10:24 kenneyf  
 Curve Type : Average

Compound	5.000 Level 1	10.000 Level 2	20.000 Level 3	50.000 Level 4	100.000 Level 5	200.000 Level 6	RRP	% RSD
	500.000 Level 7	1000.000 Level 8						
\$ 1,3,4-Dinitrotoluene	++++	46.40000	40.85000	39.74000	38.04000	38.34500	39.65376	8.405
	35.87333	38.62800						

TestAmerica West Sacramento

INITIAL CALIBRATION DATA

Start Cal Date : 09-JUN-2009 21:56  
 End Cal Date : 10-JUN-2009 04:20  
 Quant Method : ESTD  
 Origin : Disabled  
 Target Version : 4.14  
 Integrator : HP Genie  
 Method file : \\Terastation\share\GCdata\LC9.I\06092009.B\8330CNAB.M\I  
 Last Edit : 10-Jun-2009 10:25 kenneyf  
 Curve Type : Average

Calibration File Names:

Level 2: \\Terastation\share\GCdata\GCdata\LC9.I\06092009.B\C-000005.d\C-  
 Level 3: \\Terastation\share\GCdata\GCdata\LC9.I\06092009.B\C-000006.d\C-  
 Level 4: \\Terastation\share\GCdata\GCdata\LC9.I\06092009.B\C-000007.d\C-  
 Level 5: \\Terastation\share\GCdata\GCdata\LC9.I\06092009.B\C-000008.d\C-  
 Level 6: \\Terastation\share\GCdata\GCdata\LC9.I\06092009.B\C-000009.d\C-  
 Level 7: \\Terastation\share\GCdata\GCdata\LC9.I\06092009.B\C-000010.d\C-  
 Level 8: \\Terastation\share\GCdata\GCdata\LC9.I\06092009.B\C-000011.d\C-

Compound	10.000	20.000	50.000	100.000	200.000	500.000	RRF	% RSD
	Level 2	Level 3	Level 4	Level 5	Level 6	Level 7		
2 HMX	+++++	+++++	+++++	+++++	+++++	+++++	+++++	+++++
3 RDX	+++++	+++++	+++++	+++++	+++++	+++++	+++++	+++++
4 EGDN	+++++	39.00000	42.36000	44.25000	44.94000	40.36600	42.29700	5.363
5 Picric ACID	+++++	+++++	+++++	+++++	+++++	+++++	+++++	+++++
6 1,3,5-Trinitrobenzene	+++++	+++++	+++++	+++++	+++++	+++++	+++++	+++++
7 1,3-Dinitrobenzene	+++++	+++++	+++++	+++++	+++++	+++++	+++++	+++++
8 3,5-Dinitroaniline	+++++	+++++	+++++	+++++	+++++	+++++	+++++	+++++
9 TETRYL	+++++	+++++	+++++	+++++	+++++	+++++	+++++	+++++

TestAmerica West Sacramento

INITIAL CALIBRATION DATA

Start Cal Date : 09-JUN-2009 21:56  
 End Cal Date : 10-JUN-2009 04:20  
 Quant Method : ESTD  
 Origin : Disabled  
 Target Version : 4.14  
 Integrator : HP Genie  
 Method file : \\Terastation\share\GCdata\LC9.I\06092009.B\8330CNAB.M\i  
 Last Edit : 10-Jun-2009 10:25 kenneyf  
 Curve Type : Average

Compound	10 000 Level 2	20 000 Level 3	50.000 Level 4	100.000 Level 5	200.000 Level 6	500.000 Level 7	RRF	% RSD
10 Nitrobenzene	+++++	+++++	+++++	+++++	+++++	+++++	+++++	+++++
11 Nitroglycerin	57.36100	55.25000	55.16000	56.11000	57.02000	56.37800	56.23650	1.571
12 2,4,6-Trinitrotoluene	+++++	+++++	+++++	+++++	+++++	+++++	+++++	+++++
13 4-AM-2,6-DNT	+++++	+++++	+++++	+++++	+++++	+++++	+++++	+++++
14 2-AM-4,6-DNT	+++++	+++++	+++++	+++++	+++++	+++++	+++++	+++++
15 2,6-Dinitrotoluene	+++++	+++++	+++++	+++++	+++++	+++++	+++++	+++++
16 2,4-Dinitrotoluene	+++++	+++++	+++++	+++++	+++++	+++++	+++++	+++++
17 2-Nitrotoluene	+++++	+++++	+++++	+++++	+++++	+++++	+++++	+++++
18 4-Nitrotoluene	+++++	+++++	+++++	+++++	+++++	+++++	+++++	+++++
19 3-Nitrotoluene	+++++	+++++	+++++	+++++	+++++	+++++	+++++	+++++
20 PSTN	105	104	101	102	104	104	103	1.208



TestAmerica West Sacramento

INITIAL CALIBRATION DATA

Start Cal Date : 09-JUN-2009 21:56  
 End Cal Date : 10-JUN-2009 04:20  
 Quant Method : ESTD  
 Origin : Disabled  
 Target Version : 4.14  
 Integrator : HP Genie  
 Method file : \\Terastation\share\GCdata\LC9.I\06092009.B\8330CNAB.M\  
 Last Edit : 10-Jun-2009 10:25 kenneyf  
 Curve Type : Average

Compound	10.000 Level 2	20.000 Level 3	50.000 Level 4	100.000 Level 5	200.000 Level 6	500.000 Level 7	RRF	% RSD
	1000.000 Level 8							
\$ 1,3,4-Dinitrotoluene	+++++	+++++	+++++	+++++	+++++	+++++	+++++	+++++

Chromatography Summary

Injection Date: 6/10/2009 6:29 Operator: fhk  
 DataFile: LC9 I06092009.B\C-000013.D Vial Num: 20  
 Instrument ID: LC9

Method 8330 Target Analyte Results

Sample: **ICV\_6 08GCSV0397**  
**200/500/200/200/200**

Method File: LC9 I06092009 B\8330CNAB.M  
 Start Cal Date: 6/9/2009 20:53 End Cal Date: 6/10/2009 4:20

Matrix: NONE SubList: CAL.sub SpikeList:  
 Samp. Info: ICV\_6 08GCSV0397 200/500/200/200/200:2  
 Misc. Info: ; 6, . . : 3: CAL.sub ; 0. 1

Dilution Factor	Extract Volume	Sample Volume	Sample Weight
1X	0 mL	0 mL	0 g

Signal 1 UV 250-265								Signal 2 UV 358-205								
Compound Name	RT	Response	PPB	Spike Level	%D	Result	Flag	RT	Response	PPB	Spike Level	%D	Result	Flag	Limits(%)	Flag
3,4-Dinitrotoluene	31.18	259	6.5320<	200	-91%	Fails					200	-100%	Fails		(±15)	45
HMX	35.11	8323	202.0000<	200	1%	Acceptable					200	-100%	Fails		(±15)	45
RDX	25.60	8556	201.0000<	200	1%	Acceptable					200	-100%	Fails		(±15)	45
Picric ACID				500	-100%	Fails					500	-100%	Fails		(±15)	
1,3,5-Trinitrobenzene	21.78	13318	198.8000<	200	-1%	Acceptable					200	-100%	Fails		(±15)	45
1,3-Dinitrobenzene	19.50	17403	197.6000<	200	-1%	Acceptable					200	-100%	Fails		(±15)	45
TETRYL	40.80	16134	185.4000<	200	-7%	Acceptable					200	-100%	Fails		(±15)	45
Nitrobenzene	16.31	7967	201.2000<	200	1%	Acceptable					200	-100%	Fails		(±15)	45
2,4,6-Trinitrotoluene	32.92	14562	180.0000<	200	-10%	Acceptable					200	-100%	Fails		(±15)	45
4-AM-2,6-DNT	29.55	12516	187.0000<	200	-7%	Acceptable					200	-100%	Fails		(±15)	45
2-AM-4,6-DNT	29.21	14216	191.4000<	200	-4%	Acceptable					200	-100%	Fails		(±15)	45
2,6-Dinitrotoluene	27.26	9755	194.9000<	200	-3%	Acceptable					200	-100%	Fails		(±15)	45
2,4-Dinitrotoluene	26.42	15486	195.0000<	200	-3%	Acceptable					200	-100%	Fails		(±15)	45
2-Nitrotoluene	22.83	9684	401.4000<	400	0%	Acceptable					400	-100%	Fails		(±15)	45
4-Nitrotoluene				400	-100%	Fails					400	-100%	Fails		(±15)	
3-Nitrotoluene	23.35	6108	194.5000<	200	-3%	Acceptable					200	-100%	Fails		(±15)	45
Nitroglycerin				200	-100%	Fails		38.67	11785	209.6000<	200	5%	Acceptable		(±15)	45
PETN				200	-100%	Fails		49.13	19064	134.7000<	200	-8%	Acceptable		(±15)	45
3,5-Dinitroaniline	25.06	13705	204.2000<	200	2%	Acceptable					200	-100%	Fails		(±15)	45
EGDN				200	-100%	Fails		18.01	9041	213.8000<	200	7%	Acceptable		(±15)	45

Notes: M = Manually Integrated Signals Differ by More Than 40%  
 D = Operator Disabled Result Signals Differ by More Than 50%  
 O = Over Calibration Range



Chromatography Summary

Injection Date: 7/10/2009 19:32 Operator: fhk  
 DataFile: LC9 107102009A BIC-000003.D Vial Num: 2  
 Instrument ID: LC9

Method 8330 Target Analyte Results

Sample : STD\_5 09GCSV0238 .2K/1/1/1/1

Method File: LC9 107102009A BIC330CNAB M  
 Start Cal Date: 6/9/2009 20:53 End Cal Date: 6/10/2009 4:20

Matrix: NONE SubList: CAL.sub SpikeList:  
 Samp. Info: STD\_5 09GCSV0238 2K/1/1/1/1;2  
 Misc. Info: ; 5. ; , , 3; CAL.sub, , 0, 1. 0

Dilution Factor	Extract Volume	Sample Volume	Sample Weight
1X	0 mL	0 mL	0 g

Signal 1 UV 250-265								Signal 2 UV 358-205								
Compound Name	RT	Response	PPB	Spike Level	%D	Result	Flag	RT	Response	PPB	Spike Level	%D	Result	Flag	Limits(%)	Flag
3,4-Dinitrotoluene	30.47	4394	110.8000<	100	11%	Acceptable					100	-100%	Fails		(±15)	45
HMX	34.51	3416	82.8900<	100	-17%	Fails	<i>low - Not Needed</i>				100	-100%	Fails		(±15)	45
RDX	24.91	4222	99.2000<	100	-1%	Acceptable					100	-100%	Fails		(±15)	45
<b>Picric ACID</b>				<b>200</b>	<b>-100%</b>	<b>Fails</b>					<b>200</b>	<b>-100%</b>	<b>Fails</b>		<b>(±15)</b>	
1,3,5-Trinitrobenzene	21.35	6799	101.5000<	100	2%	Acceptable					100	-100%	Fails		(±15)	45
1,3-Dinitrobenzene	18.92	8855	100.6000<	100	1%	Acceptable					100	-100%	Fails		(±15)	45
TETRYL	40.33	8883	102.1000<	100	2%	Acceptable					100	-100%	Fails		(±15)	45
Nitrobenzene	16.20	4158	105.0000<	100	5%	Acceptable					100	-100%	Fails		(±15)	45
2,4,6-Trinitrotoluene	32.45	7793	96.3300<	100	-4%	Acceptable					100	-100%	Fails		(±15)	45
4-AM-2,6-DNT	28.82	6879	102.8000<	100	3%	Acceptable					100	-100%	Fails		(±15)	45
2-AM-4,6-DNT	28.50	7694	103.6000<	100	4%	Acceptable					100	-100%	Fails		(±15)	45
2,6-Dinitrotoluene	26.55	5096	101.8000<	100	2%	Acceptable					100	-100%	Fails		(±15)	45
2,4-Dinitrotoluene	25.74	8053	101.4000<	100	1%	Acceptable					100	-100%	Fails		(±15)	45
2-Nitrotoluene	22.07	5149	213.4000<	200	7%	Acceptable					200	-100%	Fails		(±15)	45
<b>4-Nitrotoluene</b>				<b>200</b>	<b>-100%</b>	<b>Fails</b>					<b>200</b>	<b>-100%</b>	<b>Fails</b>		<b>(±15)</b>	
3-Nitrotoluene	22.60	3149	100.3000<	100	0%	Acceptable					100	-100%	Fails		(±15)	45
Nitroglycerin				100	-100%	Fails		38.17	5661	100.7000<	100	1%	Acceptable		(±15)	45
PE'IN				100	-100%	Fails		48.81	10839	104.9000<	100	5%	Acceptable		(±15)	45
3,5-Dinitroaniline	24.39	6825	101.7000<	100	2%	Acceptable					100	-100%	Fails		(±15)	45
EGDN				100	-100%	Fails		17.61	4413	104.3000	100	4%	Acceptable		(±15)	45

Notes: M = Manually Integrated Signals Differ by More Than 40%  
 D = Operator Disabled Result Signals Differ by More Than 50%  
 O = Over Calibration Range

Chromatography Summary

Injection Date: 7/11/2009 5:54 Operator: flk  
 Data File: LC9 \07102009A\BIC-000013.D Vial Num: 3  
 Instrument ID: LC9

Method 8330 Target Analyte Results

Sample: **STD\_6\_09GCSV0054.SK/2/2/2/2**

Method File: LC9 \07102009A\B8330CNAB.M  
 Start Cal Date: 6/9/2009 20:53 End Cal Date: 6/10/2009 4:20

Matrix: NONE SubList: CAL.sub SpikeList:  
 Samp. Info: STD\_6\_09GCSV0054.SK/2/2/2/2  
 Misc. Info: ,6;;,3,CAL.sub;;0;1,0

Dilution Factor	Extract Volume	Sample Volume	Sample Weight
1X	0 mL	0 mL	0 g

Signal 1 UV 250-265								Signal 2 UV 358-205								
Compound Name	RT	Response	PPB	Spike Level	%D	Result	Flag	RT	Response	PPB	Spike Level	%D	Result	Flag	Limits(%)	Flag
3,4-Dinitrotoluene	30.69	8487	214.0000<	200	7%	Acceptable					200	-100%	Fails		(±15)	45
HMX	34.67	7594	184.3000<	200	-8%	Acceptable					200	-100%	Fails		(±15)	45
RDX	25.16	8673	203.8000<	200	2%	Acceptable					200	-100%	Fails		(±15)	45
<b>Picric ACID</b>				<b>500</b>	<b>-100%</b>	<b>Fails</b>					<b>500</b>	<b>-100%</b>	<b>Fails</b>		<b>(±15)</b>	
1,3,5-Trinitrobenzene	21.54	13992	208.9000<	200	4%	Acceptable					200	-100%	Fails		(±15)	45
1,3-Dinitrobenzene	19.18	18240	207.1000<	200	4%	Acceptable					200	-100%	Fails		(±15)	45
TETRYL	40.38	18246	209.7000<	200	5%	Acceptable					200	-100%	Fails		(±15)	45
Nitrobenzene	16.47	8291	209.4000<	200	5%	Acceptable					200	-100%	Fails		(±15)	45
2,4,6-Trinitrotoluene	32.59	16243	200.8000<	200	0%	Acceptable					200	-100%	Fails		(±15)	45
4-AM-2,6-DNT	29.08	13363	199.6000<	200	0%	Acceptable					200	-100%	Fails		(±15)	45
2-AM-4,6-DNT	28.74	14987	201.7000<	200	1%	Acceptable					200	-100%	Fails		(±15)	45
2,6-Dinitrotoluene	26.81	10167	203.1000<	200	2%	Acceptable					200	-100%	Fails		(±15)	45
2,4-Dinitrotoluene	25.99	16528	208.1000<	200	4%	Acceptable					200	-100%	Fails		(±15)	45
2-Nitrotoluene	22.38	10164	421.3000<	400	5%	Acceptable					400	-100%	Fails		(±15)	45
<b>4-Nitrotoluene</b>				<b>400</b>	<b>-100%</b>	<b>Fails</b>					<b>400</b>	<b>-100%</b>	<b>Fails</b>		<b>(±15)</b>	
3-Nitrotoluene	22.90	6323	201.4000<	200	1%	Acceptable					200	-100%	Fails		(±15)	45
<b>Nitroglycerin</b>				<b>200</b>	<b>-100%</b>	<b>Fails</b>		38.25	11803	209.9000<	200	5%	Acceptable		(±15)	45
<b>PETN</b>				<b>200</b>	<b>-100%</b>	<b>Fails</b>		48.78	21503	208.1000<	200	4%	Acceptable		(±15)	45
3,5-Dinitroaniline	24.65	13757	205.0000<	200	3%	Acceptable					200	-100%	Fails		(±15)	45
<b>EGDN</b>				<b>200</b>	<b>-100%</b>	<b>Fails</b>		17.80	9381	221.8000	200	11%	Acceptable		(±15)	45

Notes: M = Manually Integrated Signals Differ by More Than 40%  
 D = Operator Disabled Result Signals Differ by More Than 50%  
 O = Over Calibration Range



Chromatography Summary

Injection Date: 7/14/2009 15:00 Operator: flk  
 DataFile: LC9 R07142009A BAC-000002.D Vial Num: 2  
 Instrument ID: LC9

Method 8330 Target Analyte Results

Sample: STD\_5 09GCSV0238 .2K/1/1/1/1

Method File: LC9 R07142009A B8330CNAB.M  
 Start Cal Date: 6/9/2009 20:53 End Cal Date: 6/10/2009 4:20

Matrix: NONE SubList: CAL.sub SpikeList:  
 Samp. Info: STD\_5 09GCSV0238 .2K/1/1/1/1,2  
 Misc. Info: ; 5 ; ; 3, CAL.sub ; 0 ; 1 ; 0

Dilution Factor	Extract Volume	Sample Volume	Sample Weight
1X	0 mL	0 mL	0 g

Signal 1 UV 250-265								Signal 2 UV 358-205								
Compound Name	RT	Response	PPB	Spike Level	%D	Result	Flag	RT	Response	PPB	Spike Level	%D	Result	Flag	Limits(%)	Flag
3,4-Dinitrotoluene	31.08	4056	102.3000<	100	2%	Acceptable					100	-100%	Fails		(±15)	45
HMX	35.02	4112	99.7800<	100	0%	Acceptable					100	-100%	Fails		(±15)	45
RDX	25.52	4319	101.5000<	100	2%	Acceptable					100	-100%	Fails		(±15)	45
<b>Picric ACID</b>				<b>200</b>	<b>-100%</b>	<b>Fails</b>					<b>200</b>	<b>-100%</b>	<b>Fails</b>		<b>(±15)</b>	
1,3,5-Trinitrobenzene	21.79	6774	101.1000<	100	1%	Acceptable					100	-100%	Fails		(±15)	45
1,3-Dinitrobenzene	19.46	8818	100.1000<	100	0%	Acceptable					100	-100%	Fails		(±15)	45
TETRYL	40.68	9087	104.4000<	100	4%	Acceptable					100	-100%	Fails		(±15)	45
Nitrobenzene	16.75	4130	104.3000<	100	4%	Acceptable					100	-100%	Fails		(±15)	45
2,4,6-Trinitrotoluene	32.85	7977	98.6100<	100	-1%	Acceptable					100	-100%	Fails		(±15)	45
4-AM-2,6-DNT	29.46	6865	102.5000<	100	3%	Acceptable					100	-100%	Fails		(±15)	45
2-AM-4,6-DNT	29.11	7562	101.8000<	100	2%	Acceptable					100	-100%	Fails		(±15)	45
2,6-Dinitrotoluene	27.17	5110	102.1000<	100	2%	Acceptable					100	-100%	Fails		(±15)	45
2,4-Dinitrotoluene	26.34	8075	101.7000<	100	2%	Acceptable					100	-100%	Fails		(±15)	45
2-Nitrotoluene	22.75	5063	209.8000<	200	5%	Acceptable					200	-100%	Fails		(±15)	45
<b>4-Nitrotoluene</b>				<b>200</b>	<b>-100%</b>	<b>Fails</b>					<b>200</b>	<b>-100%</b>	<b>Fails</b>		<b>(±15)</b>	
3-Nitrotoluene	23.27	3193	101.7000<	100	2%	Acceptable					100	-100%	Fails		(±15)	45
<b>Nitroglycerin</b>				<b>100</b>	<b>-100%</b>	<b>Fails</b>		38.56	5792	103.0000<	100	3%	Acceptable		(±15)	45
<b>PETN</b>				<b>100</b>	<b>-100%</b>	<b>Fails</b>		49.00	10860	105.1000<	100	5%	Acceptable		(±15)	45
3,5-Dinitroaniline	25.01	6927	103.2000<	100	3%	Acceptable					100	-100%	Fails		(±15)	45
<b>EGDN</b>				<b>100</b>	<b>-100%</b>	<b>Fails</b>		18.00	4612	109.0000	100	9%	Acceptable		(±15)	45

Notes: M = Manually Integrated Signals Differ by More Than 40%  
 D = Operator Disabled Result Signals Differ by More Than 50%  
 O = Over Calibration Range

Chromatography Summary

Injection Date: 7/14/2009 17:04 Operator: fhk  
 DataFile: LC9.R07142009A.BAC-000004.D Vial Num: 3  
 Instrument ID: LC9

Method 8330 Target Analyte Results

Sample: STD\_6 09GCSV0054 .5K/2/2/2/2

Method File: LC9.R07142009A.B8330CNAB.M  
 Start Cal Date: 6/9/2009 20:53 End Cal Date: 6/10/2009 4:20

Matrix: NONE SubList: CAL.sub SpikeList:  
 Samp. Info: STD\_6 09GCSV0054 .5K/2/2/2/2  
 Misc. Info: .6, ; ; 3, CAL.sub; ; 0; 1; 0

Dilution Factor	Extract Volume	Sample Volume	Sample Weight
1X	0 mL	0 mL	0 g

Compound Name	RT	Signal 1 UV 250-265					Signal 2 UV 358-205					Limits(%)	Flag			
		Response	PPB	Spike Level	%D	Result	RT	Response	PPB	Spike Level	%D			Result		
3,4-Dinitrotoluene	31.17	8088	204.0000<	200	2%	Acceptable						200	-100%	Fails	(±15)	45
HMX	35.10	8294	201.3000<	200	1%	Acceptable						200	-100%	Fails	(±15)	45
RDX	25.64	8841	207.7000<	200	4%	Acceptable						200	-100%	Fails	(±15)	45
<b>Picric ACID</b>				<b>500</b>	<b>-100%</b>	<b>Fails</b>						<b>500</b>	<b>-100%</b>	<b>Fails</b>	<b>(±15)</b>	
1,3,5-Trinitrobenzene	21.85	13960	208.4000<	200	4%	Acceptable						200	-100%	Fails	(±15)	45
1,3-Dinitrobenzene	19.56	18224	207.0000<	200	4%	Acceptable						200	-100%	Fails	(±15)	45
TETRYL	40.73	18387	211.3000<	200	6%	Acceptable						200	-100%	Fails	(±15)	45
Nitrobenzene	16.87	8317	210.0000<	200	5%	Acceptable						200	-100%	Fails	(±15)	45
2,4,6-Trinitrotoluene	32.91	15780	195.1000<	200	-2%	Acceptable						200	-100%	Fails	(±15)	45
4-AM-2,6-DNT	29.58	13774	205.7000<	200	3%	Acceptable						200	-100%	Fails	(±15)	45
2-AM-4,6-DNT	29.23	15080	203.0000<	200	2%	Acceptable						200	-100%	Fails	(±15)	45
2,6-Dinitrotoluene	27.29	10345	206.7000<	200	3%	Acceptable						200	-100%	Fails	(±15)	45
2,4-Dinitrotoluene	26.45	16604	209.1000<	200	5%	Acceptable						200	-100%	Fails	(±15)	45
2-Nitrotoluene	22.87	10117	419.3000<	400	5%	Acceptable						400	-100%	Fails	(±15)	45
<b>4-Nitrotoluene</b>				<b>400</b>	<b>-100%</b>	<b>Fails</b>						<b>400</b>	<b>-100%</b>	<b>Fails</b>	<b>(±15)</b>	
3-Nitrotoluene	23.38	6428	204.7000<	200	2%	Acceptable						200	-100%	Fails	(±15)	45
<b>Nitroglycerin</b>				<b>200</b>	<b>-100%</b>	<b>Fails</b>	38.61	12038	214.1000<	200	7%	Acceptable			(±15)	45
<b>PETN</b>				<b>200</b>	<b>-100%</b>	<b>Fails</b>	48.99	21884	211.8000<	200	6%	Acceptable			(±15)	45
3,5-Dinitroaniline	25.12	13891	207.0000<	200	4%	Acceptable						200	-100%	Fails	(±15)	45
<b>EGDN</b>				<b>200</b>	<b>-100%</b>	<b>Fails</b>	18.07	9241	218.5000	200	9%	Acceptable			(±15)	45

Notes M = Manually Integrated Signals Differ by More Than 40%  
 D = Operator Disabled Result Signals Differ by More Than 50%  
 O = Over Calibration Range



2324MWO1AW

TestAmerica West Sacramento

A-000006.D

Chromatography Summary

Injection Date: 7/8/2009 20:22 Operator: fsk  
Data File: LC10 107082009.BVA-000006.D Vial Num: 13  
Instrument ID: LC10

Method 8330 Target Analyte Results

Sample: LFXXW1AC 9183251 G9G010225-1  
1X

Method File: LC10 107082009.BV330AB.M  
Start Cal Date: 6/10/2009 15:33 End Cal Date: 6/11/2009 13:47

Matrix: WATER SubList: WATER.sub SpikeList:  
Samp. Info: LFXXW1AC 9183251 G9G010225-1 1X;0;  
Misc. Info: .:1021 57;:20:1;WATER sub;:0,1;LFXXW1AC

Dilution Factor	Extract Volume	Sample Volume	Sample Weight
1X	20 mL	1021.57 mL	0 g

Signal 1 UV 250-265						Signal 2 UV 358-205							
Compound Name	RT	Diff	Response	Conc (ug/L)	Flag	RT	Diff	Response	Conc (ug/L)	Flag	MDL	RL	Flag
3,4-Dinitrotoluene	18.24	0.032	5380	2.2330<		18.24	0.029	10484	2.3740		0.0000	0.00	
HMX											0.0264	0.10	
RDX	7.88	0.032	524	0.1317<	confirmed once						0.0636	0.10	45
Picric ACID											0.1224	0.98	
1,3,5-Trinitrobenzene	10.12	-0.088	247	0.0328<							0.0303	0.10	45
1,3-Dinitrobenzene											0.0489	0.10	
TETRYL											0.0489	0.10	
Nitrobenzene											0.0489	0.10	
2,4,6-Trinitrotoluene											0.0235	0.10	
4-AM-2,6-DNT											0.0215	0.10	
2-AM-4,6-DNT											0.0979	0.20	
2,6-Dinitrotoluene											0.0489	0.10	
2,4-Dinitrotoluene											0.0489	0.10	
2-Nitrotoluene											0.0705	0.49	
4-Nitrotoluene											0.0705	0.49	
3-Nitrotoluene											0.0607	0.49	
Nitroglycenn											0.3230	0.64	
PETN											0.2937	0.64	
3,5-Dinitroaniline											0.0245	0.98	

Surrogates:	Spiked	Recovered	% Rec	Spiked	Recovered	% Rec	Limits
3,4-Dinitrotoluene	2.4472	2.2330	91	2.4472	2.3740	97	(79-111)

Notes: M = Manually Integrated Signals Differ by More Than 40%  
D = Operator Disabled Result Signals Differ by More Than 50%  
O = Over Calibration Range

Chromatography Summary

Injection Date: 7/10/2009 21:36 Operator: fhk  
 DataFile: I:\91107102009A\BAC-000005.D Vial Num: 12  
 Instrument ID: 1C9

Method 8330 Target Analyte Results

Sample: **LFXXWIAC 9183251 G9G010225-1**  
**1X**

Method File: I:\91107102009A\B8330CNAB.M  
 Start Cal Date: 6/9/2009 20:53 End Cal Date: 6/10/2009 4:20

Matrix: WATER SubList: WATER.sub SpikeList:  
 Samp. Info: LFXXWIAC 9183251 G9G010225-1 1X;0;  
 Misc. Info: ;:1021.57;;20;1;WATER.sub;;0;1.LFXXWIAC

Dilution Factor	Extract Volume	Sample Volume	Sample Weight
1X	20 mL	1021.57 mL	0 g

Signal 1 UV 250-265						Signal 2 UV 358-205							
Compound Name	RT	Diff	Response	Conc (ug/L)	Flag	RT	Diff	Response	Conc (ug/L)	Flag	MDL	RL	Flag
3,4-Dinitrotoluene	30.52	-0.138	4974	2.4560<							0.0000	0.00	45
HMX											0.0264	0.10	
RDX	24.99	-0.140	269	0.1237<	con. turned						0.0636	0.10	45
Picric ACID											0.1224	0.96	
1,3,5-Trinitrobenzene											0.0303	0.10	
1,3-Dinitrobenzene											0.0489	0.10	
TETRYL											0.0489	0.10	
Nitrobenzene											0.0489	0.10	
2,4,6-Trinitrotoluene											0.0235	0.10	
4-AM-2,6-DNT											0.0215	0.10	
2-AM-4,6-DNT											0.0979	0.19	
2,6-Dinitrotoluene											0.0489	0.10	
2,4-Dinitrotoluene											0.0489	0.10	
2-Nitrotoluene											0.0705	0.48	
4-Nitrotoluene											0.0705	0.48	
3-Nitrotoluene											0.0607	0.48	
Nitroglycerin											0.3230	0.62	
PETN											0.2937	0.62	
3,5-Dinitroaniline											0.0245	0.96	

Surrogates:	Spiked	Recovered	% Rec	Spiked	Recovered	% Rec	Limits
3,4-Dinitrotoluene	2.4472	2.4560	100	2.4472	0		(48 143)

Notes M = Manually Integrated Signals Differ by More Than 40%  
 D = Operator Disabled Result Signals Differ by More Than 50%  
 O = Over Calibration Range

Chromatography Summary

Injection Date: 7/8/2009 21:12 Operator: fhk  
Data File: LC10 R07082009 BAA-000007.D Vial Num: 14  
Instrument ID: LC10

Method 8330 Target Analyte Results

Sample: LFXX61AC 9183251 G9G010225-2 1X

Method File: LC10 R07082009 B8330AB.M  
Start Cal Date: 6/10/2009 15:33 End Cal Date: 6/11/2009 13:47

Matrix: WATER SubList: WATER.sub SpikeList:  
Samp. Info: LFXX61AC 9183251 G9G010225-2 1X.0,  
Misc. Info: ;1022.4.;20.1;WATER sub.;0.1:LFXX61AC

Dilution Factor	Extract Volume	Sample Volume	Sample Weight
1X	10 mL	1022.4 mL	0 g

Signal 1 UV 250-265						Signal 2 UV 358-205							
Compound Name	RT	Diff	Response	Conc (ug/L)	Flag	RT	Diff	Response	Conc (ug/L)	Flag	MDL	RL	Flag
3,4-Dinitrotoluene	18.22	0.008	5488	2.2760<		18.22	0.008	10601	2.3990		0.0000	0.00	
HMX											0.0264	0.10	
RDX											0.0616	0.10	
Picric ACID											0.1223	0.98	
1,3,5-Trinitrobenzene	10.11	-0.089	252	0.033<<							0.0303	0.10	45
1,3-Dinitrobenzene											0.0489	0.10	
TETRYL											0.0489	0.10	
Nitrobenzene											0.0489	0.10	
2,4,6-Trinitrotoluene											0.0235	0.10	
4-AM-2,6-DNT											0.0215	0.10	
2-AM-4,6-DNT											0.0978	0.20	
2,6-Dinitrotoluene											0.0489	0.10	
2,4-Dinitrotoluene											0.0489	0.10	
2-Nitrotoluene											0.0704	0.49	
4-Nitrotoluene											0.0704	0.49	
3-Nitrotoluene											0.0606	0.49	
Nitroglycerin											0.3228	0.64	
PETN											0.2934	0.64	
3,5-Dinitroaniline											0.0245	0.98	

Surrogates:	Spiked	Recovered	% Rec	Spiked	Recovered	% Rec	Limits
3,4-Dinitrotoluene	2.4452	2.2760	93	2.4452	2.3990	98	(79-111)

Notes: M = Manually Integrated Signals Differ by More Than 40%  
D = Operator Disabled Result Signals Differ by More Than 50%  
O = Over Calibration Range

*not confirmed*

23L4MW460W

A-000008.D

Chromatography Summary

Injection Date: 7/8/2009 22:03 Operator: fhk  
 Data File: I.C10 N07082009 BVA-000008.D Vial Num: 15  
 Instrument ID: I.C10

Method 8330 Target Analyte Results

Sample: LFX111AC 9183251 G9G010225-3 1X

Method File: I.C10 N07082009 B8330AB.M  
 Start Cal Date: 6/10/2009 15:33 End Cal Date: 6/11/2009 13:47

Dilution Factor	Extract Volume	Sample Volume	Sample Weight
1X	20 mL	1024.15 mL	0 g

Matrix: WATER SubList: WATER sub SpikeList:  
 Samp. Info: LFX111AC 9183251 G9G010225-3 1X:0  
 Misc. Info: ..1024 15.;20.1.WATER sub.,0.1.LFX111AC

Signal 1 UV 250-265						Signal 2 UV 358-205							
Compound Name	RT	Diff	Response	Conc (ug/L)	Flag	RT	Diff	Response	Conc (ug/L)	Flag	MDL	RL	Flag
3,4-Dinitrotoluene	18.22	0.013	5492	2.2740<		18.23	0.017	10608	2.3960		0.0000	0.00	
HMX	5.30	0.007	178	0.0288<							0.0264	0.10	45
RDX	7.87	0.013	16452	4.1250<							0.0635	0.10	45
Picric ACID											0.1221	0.98	
1,3,5-Trinitrobenzene	10.36	0.160	1266	0.1675<							0.0303	0.10	45
1,3-Dinitrobenzene											0.0488	0.10	
TETRYL											0.0488	0.10	
Nitrobenzene											0.0488	0.10	
2,4,6-Trinitrotoluene											0.0234	0.10	
4-AM-2,6-DNT											0.0215	0.10	
2-AM-4,6-DNT											0.0976	0.20	
2,6-Dinitrotoluene											0.0488	0.10	
2,4-Dinitrotoluene											0.0488	0.10	
2-Nitrotoluene											0.0703	0.49	
4-Nitrotoluene											0.0703	0.49	
3-Nitrotoluene											0.0605	0.49	
Nitroglycerin											0.3222	0.63	
PETN											0.2929	0.63	
3,5-Dinitroaniline											0.0244	0.98	

*contributed on cu*  
*not contributed on cu*  
 7/14/09

Surrogates:	Spiked	Recovered	% Rec	Spiked	Recovered	% Rec	Limits
3,4-Dinitrotoluene	2.4410	2.2740	93	2.4410	2.3960	98	(79-111)

Notes: M = Manually Integrated Signals Differ by More Than 40%  
 D = Operator Disabled Result Signals Differ by More Than 50%  
 O = Over Calibration Range

Chromatography Summary

Injection Date: 7/10/2009 22:39 Operator: fbk  
 DataFile: LC9 107102009A BIC-000006.D Vial Num: 13  
 Instrument ID: LC9

Method 8330 Target Analyte Results

Sample: LFX111AC 9183251 G9G010225-3 1X Method File: LC9 107102009A B\8330CNAB.M  
 Start Cal Date: 6/9/2009 20:53 End Cal Date: 6/10/2009 4:20

Dilution Factor	Extract Volume	Sample Volume	Sample Weight
1X	20 mL	1024.15 mL	0 g

Matrix: WATER SubList: WATER sub SpikeList:  
 Samp. Info: LFX111AC 9183251 G9G010225-3 1X;0;  
 Misc. Info: ;;1024 15,,20;1,WATER sub,,0;1;LFX111AC

Signal 1 UV 250-265						Signal 2 UV 358-205							
Compound Name	RT	Diff	Response	Conc (ug/L)	Flag	RT	Diff	Response	Conc (ug/L)	Flag	MDL	RL	Flag
3,4-Dinitrotoluene	30.74	0.087	4980	2.4520<							0.0000	0.00	45
HMX	34.72	0.068	89	0.0422<							0.0264	0.10	45
RDX	25.26	0.125	8280	3.7990<	- confirmed						0.0635	0.10	45
Picric ACID											0.1221	0.95	
1,3,5-Trinitrobenzene					- Not confirmed MD						0.0303	0.10	
1,3-Dinitrobenzene											0.0488	0.10	
TETRYL											0.0488	0.10	
Nitrobenzene											0.0488	0.10	
2,4,6-Trinitrotoluene											0.0234	0.10	
4-AM-2,6-DNT											0.0215	0.10	
2-AM-4,6-DNT											0.0976	0.19	
2,6-Dinitrotoluene											0.0488	0.10	
2,4-Dinitrotoluene											0.0488	0.10	
2-Nitrotoluene											0.0703	0.48	
4-Nitrotoluene											0.0703	0.48	
3-Nitrotoluene											0.0605	0.48	
Nitroglycerin											0.3222	0.62	
PETN											0.2929	0.62	
3,5-Dinitroaniline											0.0244	0.95	

*MD*  
 7/14/09

Surrogate(s)	Spiked	Recovered	% Rec	Spiked	Recovered	% Rec	Limits
3,4-Dinitrotoluene	2.4410	2.4520	100	2.4410		0	(48-143)

Notes: M = Manually Integrated Signals Differ by More Than 40%  
 D = Operator Disabled Result Signals Differ by More Than 50%  
 O = Over Calibration Range

2324 MWO 3BW

TestAmerica West Sacramento

A-000009.D

Chromatography Summary

Injection Date: 7/8/2009 22:53 Operator: flik  
DataFile: LC10 1107082009 BVA-000009.D Vial Num: 16  
Instrument ID: LC10

Method 8330 Target Analyte Results

Sample: LFX121AC 9183251 G9G010225-4 1X

Method File: LC10 1107082009 B8330AB.M  
Start Cal Date: 6/10/2009 15:33 End Cal Date: 6/11/2009 13:47

Dilution Factor	Extract Volume	Sample Volume	Sample Weight
1X	20 mL	1019.2 mL	0 g

Matrix: WATER SubList: WATER.sub SpikeList:  
Samp. Info: LFX121AC 9183251 G9G010225-4 1X,0  
Misc. Info: ;,1019 2.,20,1,WATER sub,,0 1;LFX121AC

Signal 1 UV 250-265					Signal 2 UV 358-205								
Compound Name	RT	Diff	Response	Conc (ug/L)	Flag	RT	Diff	Response	Conc (ug/L)	Flag	MDL	RL	Flag
3,4-Dinitrotoluene	18.24	0.030	5457	2.270<		18.24	0.033	10513	2.3860		0.0000	0.00	
HMX	5.30	0.006	169	0.027<							0.0265	0.10	45
RDX	7.87	0.020	16232	4.0890<							0.0638	0.10	45
Picric ACID											0.1226	0.98	
1,3,5-Trinitrobenzene	10.39	0.183	1769	0.2352<							0.0304	0.10	45
1,3-Dinitrobenzene											0.0491	0.10	
TETRYL											0.0491	0.10	
Nitrobenzene											0.0491	0.10	
2,4,6-Trinitrotoluene											0.0235	0.10	
4-AM-2,6-DNT											0.0216	0.10	
2-AM-4,6-DNT											0.0981	0.20	
2,6-Dinitrotoluene											0.0491	0.10	
2,4-Dinitrotoluene											0.0491	0.10	
2-Nitrotoluene											0.0706	0.49	
4-Nitrotoluene											0.0706	0.49	
3-Nitrotoluene											0.0608	0.49	
Nitroglycerin											0.3238	0.64	
PETN											0.2943	0.64	
3,5-Dinitroaniline											0.0245	0.98	

*contaminated cu*  
*not contaminated cu*  
*gfk*  
*7/14/09*

Surrogates:	Spiked	Recovered	% Rec	Spiked	Recovered	% Rec	Limits
3,4-Dinitrotoluene	2.4529	2.2700	93	2.4529	2.3860	97	(79-111)

Notes: M = Manually Integrated Signals Differ by More Than 40%  
D = Operator Disabled Result Signals Differ by More Than 50%  
O = Over Calibration Range

Chromatography Summary

Injection Date: 7/10/2009 23:41 Operator: ftk  
 DataFile: LC9 I07102009A BAC-000007.D Vial Num: 14  
 Instrument ID: LC9

Method 8330 Target Analyte Results

Sample: LFX121AC 9183251 G9G010225-4 1X

Method File: LC9 I07102009A B8330CNAB.M  
 Start Cal Date: 6/9/2009 20:53 End Cal Date: 6/10/2009 4:20

Dilution Factor	Extract Volume	Sample Volume	Sample Weight
1X	20 mL	1019.2 mL	0 g

Matrix: WATER SubList: WATER sub SpikeList:  
 Samp. Info: LFX121AC 9183251 G9G010225-4 1X,0,  
 Misc. Info: ;,1019 2,;20;1;WATER.sub;0;1;LFX121AC

Signal 1 UV 250-265						Signal 2 UV 358-205							
Compound Name	RT	Diff	Response	Conc (ug/L)	Flag	RT	Diff	Response	Conc (ug/L)	Flag	MDL	RL	Flag
3,4-Dinitrotoluene	30.40	-0.259	5098	2.5230<	M						0.0000	0.00	45
<u>HMX</u>											0.0265	0.10	
RDX	24.85	-0.286	8046	3.7100<	M						0.0638	0.10	45
Picric ACID											0.1226	0.96	
1,3,5-Trinitrobenzene											0.0304	0.10	
1,3-Dinitrobenzene	18.86	-0.292	1390	0.3098<	M						0.0491	0.10	45
TETRYL											0.0491	0.10	
Nitrobenzene											0.0491	0.10	
2,4,6-Trinitrotoluene											0.0235	0.10	
1-AM-2,6-DNT	28.72	-0.320	96	0.0281<	M						0.0216	0.10	45
2-AM-4,6-DNT											0.0981	0.19	
2,6-Dinitrotoluene											0.0491	0.10	
2,4-Dinitrotoluene											0.0491	0.10	
2-Nitrotoluene											0.0706	0.48	
4-Nitrotoluene											0.0706	0.48	
3-Nitrotoluene											0.0608	0.48	
Nitrolycerin											0.3238	0.63	
PETN											0.2943	0.63	
3,5-Dinitroaniline											0.0245	0.96	

*Continued*

*Not continued (M)*

*7/14/09*

*reprocess*

Surrogates:	Spiked	Recovered	% Rec	Spiked	Recovered	% Rec	Limits
3,4-Dinitrotoluene	2.4529	2.5230	103	2.4529		0	(48 143)

Notes M :: Manually Integrated Signals Differ by More Than 40%  
 D :: Operator Disabled Result Signals Differ by More Than 50%  
 O :: Over Calibration Range

2324 MW 08AW

TestAmerica West Sacramento

A-000010.D

Chromatography Summary

Injection Date: 7/8/2009 23:44 Operator: fhk  
Data File: LC10 B07082009 BVA-000010.D Vial Num: 17  
Instrument ID: LC10

Method 8330 Target Analyte Results

Sample: LFX151AC 9183251 G9G010225-5 IX

Method File: LC10 B07082009.B8330AB.M  
Start Cal Date: 6/10/2009 15:33 End Cal Date: 6/11/2009 13:47

Matrix: WATER SubList: WATER sub SpikeList:  
Samp. Info: LFX151AC 9183251 G9G010225-5 IX,0,  
Misc. Info: ;1021.92;20:1.WATER sub,,0.1.LFX151AC

Dilution Factor	Extract Volume	Sample Volume	Sample Weight
1X	20 mL	1021.92 mL	0 g

Signal 1 UV 250-265						Signal 2 UV 358-205							
Compound Name	RT	Diff	Response	Conc (ug/L)	Flag	RT	Diff	Response	Conc (ug/L)	Flag	MDL	RL	Flag
3,4-Dinitrotoluene	18.23	0.021	5597	2.3220<		18.24	0.025	10764	2.4370		0.0000	0.00	
HMX	5.30	0.005	2345	0.3798<							0.0264	0.10	45
RDX	7.87	0.018	37290	9.3690<							0.0636	0.10	45
Picric ACID											0.1223	0.98	
1,3,5-Trinitrobenzene	10.38	0.171	1655	0.2194<							0.0303	0.10	45
1,3-Dinitrobenzene											0.0489	0.10	
TETRYL											0.0489	0.10	
Nitrobenzene											0.0489	0.10	
2,4,6-Trinitrotoluene											0.0235	0.10	
4-AM-2,6-DNT											0.0215	0.10	
2-AM-4,6-DNT											0.0979	0.20	
2,6-Dinitrotoluene											0.0489	0.10	
2,4-Dinitrotoluene											0.0489	0.10	
2-Nitrotoluene											0.0705	0.49	
4-Nitrotoluene											0.0705	0.49	
3-Nitrotoluene											0.0607	0.49	
Nitroglycerin											0.3229	0.64	
PETN											0.2936	0.64	
3,5-Dinitroaniline											0.0245	0.98	

*Conc. increased in air*  
*Not confirmed on air*  
*gmk 7/14/09*

Surrogates:	Spiked	Recovered	% Rec	Spiked	Recovered	% Rec	Limits
3,4-Dinitrotoluene	2.4464	2.3220	95	2.4464	2.4370	100	(79-111)

Notes: M = Manually Integrated Signals Differ by More Than 40%  
D = Operator Disabled Result Signals Differ by More Than 50%  
O = Over Calibration Range



Chromatography Summary

Injection Date: 7/14/2009 16:02 Operator: fhk  
 DataFile: LC9\107142009A BIC-000003.D Vial Num: 11  
 Instrument ID: LC9

Method 8330 Target Analyte Results

Sample: LFX151AC 9183251 G9G010225-5 1X

Method File: LC9\107142009A B18330CNAB.M  
 Start Cal Date: 6/9/2009 20:53 End Cal Date: 6/10/2009 4:20

Matrix: WATER SubList: WATER sub SpikeList:  
 Samp. Info: LFX151AC 9183251 G9G010225-5 1X;0;  
 Misc. Info: ..1021 92;;20;1;WATER.sub;0;1;LFX151AC

Dilution Factor	Extract Volume	Sample Volume	Sample Weight
1X	20 mL	1021.92 mL	0 g

Compound Name	Zorbax Cyano(250nm)					Zorbax Cyano(250nm-205nm)					MDL	RL	Flag
	RT	Diff	Response	Conc (ug/L)	Flag	RT	Diff	Response	Conc (ug/L)	Flag			
3,4-Dinitrotoluene	31.20	0.030	4881	2.4090<							0.0000	0.00	45
HMX	38.09	-0.006	823	0.3908<							0.0264	0.10	45
RDX	25.67	0.032	18997	8.7360<							0.0636	0.10	45
Picric ACID											0.1223	0.96	
1,3,5-Trinitrobenzene											0.0303	0.10	
1,3-Dinitrobenzene											0.0489	0.10	
TETRYL											0.0489	0.10	
Nitrobenzene											0.0489	0.10	
2,4,6-Trinitrotoluene											0.0235	0.10	
4-AM-2,6-DNT	29.66	0.075	84	0.0246<							0.0215	0.10	45
2-AM-4,6-DNT											0.0979	0.19	
2,6-Dinitrotoluene											0.0489	0.10	
2,4-Dinitrotoluene											0.0489	0.10	
2-Nitrotoluene											0.0705	0.48	
4-Nitrotoluene											0.0705	0.48	
3-Nitrotoluene											0.0607	0.48	
Nitroglycerin											0.3229	0.62	
PETN											0.2936	0.62	
3,5-Dinitroaniline											0.0245	0.96	

*not confirmed*  
*confirmed*  
 JUN 5/15/09

Surrogates:	Spiked	Recovered	% Rec	Spiked	Recovered	% Rec	Limits
3,4-Dinitrotoluene	2.4464	2.4090	98	2.4464	0	0	(48-143)

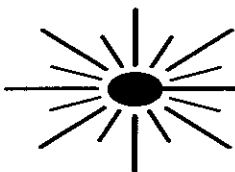
Notes M = Manually Integrated Signals Differ by More Than 40%  
 D = Operator Disabled Result Signals Differ by More Than 50%  
 O = Over Calibration Range

**Data Validation Report**

**Michael Baker Jr., Inc**

**Camp Bonneville**

**SDG#: PSF0979**

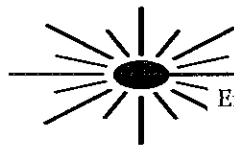


**ECT.CON INC.**

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Environmental and Computer  
Technology Consultants

3531 Fox Chase Drive  
Imperial, PA 15126  
(724) 695-8042  
FAX (724) 695-2698  
e-mail: [ectconinc@comcast.net](mailto:ectconinc@comcast.net)



## Data Validation Report

SDG#	PSF0979
Validation Report Date	August 17, 2009
Validation Guidance	USEPA CLP National Functional Guidelines for Data Review
Client Name	Michael Baker Jr., Inc.
Project Name	Camp Bonneville
Laboratory	TestAmerica
Analytical Parameters	Explosives
Analytical Method	SW-846 8330

### Samples/Matrix:

Date Sampled	Location	Portland ID	Sacramento ID	Explosives	Matrix
06/29/09	23L4MW04AW	PSF0979-01	LF3H4	X	Aqueous
06/29/09	23L4MW02AW	PSF0979-02	LF3H5	X	Aqueous
06/29/09	23L4MW02BW	PSF0979-03	LF3H6	X	Aqueous

Analytical data in this report were screened to determine analytical limitations of the data based on specific quality control criteria. This screening assumes analytical results are correct as reported and merely provides an interpretation of the reported quality control results. Laboratory calculations have been verified as part of this validation. Specific findings on analytical limitations are presented in this report. Annotated Form 1s or spreadsheets for samples reviewed are included after the Data Assessment Findings. Form 1s for the MS/MSD samples and spreadsheets are not annotated.

### SUMMARY

The sample set for the Camp Bonneville site consists of three aqueous field samples. These samples were analyzed for explosives as provided in the Table 1. The findings presented in this review of the analytical data assume that the information presented by the analytical laboratory is correct.

The explosives findings are based upon the assessment of the following:

- \* ● Data Completeness
- \* ● Holding Times
- \* ● Calibration (Initial and Continuing)
- \* ● Blanks
- \* ● System Monitoring Compounds (Surrogate Spikes)
- \* ● Matrix Spike/Matrix Spike Duplicates
- Laboratory Control Standard Results
- \* ● Target Compound Identification
- Compound Quantification and Reported Contract Quantitation Limits
- \* ● System Performance

\* Criteria were met for this evaluation item.

This evaluation was conducted in accordance with USEPA CLP National Functional Guidelines for Organic Data Review and the analytical method. Findings from this evaluation should be considered when using the analytical data. This report presents a summary of the data qualifications based on the review of the aforementioned evaluation criteria. This is followed by annotated Form 1s/ spreadsheets. Finally, the worksheets used to perform the evaluation are provided.

## **FINDINGS**

### **1. Laboratory Control Sample Results**

Recovery of HMX (116%) and RDX (124%) exceeded the laboratory's upper quality control limits. In the following samples, positive results for HMX and RDX were qualified as estimated "J".

23L4MW04AW          23L4MW02AW          23L4MW02BW

### **2. Compound Quantitation**

Positive results less than the reporting limit were qualified as estimated "J" due to the level of uncertainty.

## **NOTES**

### **Matrix Spike/Matrix Spike Duplicate Results**

A MS/MSD was not analyzed with this SDG. A laboratory control sample was analyzed. Data were not qualified on this basis.

### **Laboratory Control Sample Results**

Recovery of several compounds exceeded the laboratory's upper quality control limits. The affected compounds were not detected in the associated samples. Data were not qualified on this basis.

### **Compound Quantitation**

Sample 23L4MW02BW was analyzed and reported at a 5X dilution factor due to the presence of RDX above the linear calibration range of the instrument. This accounts for the elevated reporting limits for this sample. Data were not qualified on this basis.

### **Field Duplicate Results**

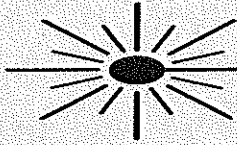
No field duplicates were included with this SDG. Data were not qualified on this basis.

  
Data Reviewer

8/17/09  
Date

### **Glossary of Data Qualifiers**

U	Not Detected.	The associated number indicates approximate sample concentration necessary to be detected.
UJ	Not Detected.	Quantitation limit may be inaccurate or imprecise.
J	Analyte Present.	Reported value may not be accurate or precise.
N	Consider Present.	Tentative identification. Special methods may be needed to confirm its presence or absence in future sampling efforts.
R	Unusable Result.	Analyte may or may not be present in the sample.
UR	Unusable Result.	Analyte may or may not be present in the sample.



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*Annotated Form 1's*  
*(Spreadsheet)*

TestAmerica Portland

Client Sample ID: PSF0979-01

HPLC 23L4MW04AW

Lot-Sample #...: G9G030156-001    Work Order #...: LF3H41AC    Matrix.....: WATER  
 Date Sampled...: 06/29/09    Date Received...: 07/02/09  
 Prep Date.....: 07/06/09    Analysis Date...: 07/09/09  
 Prep Batch #...: 9187206    *0306*  
 Dilution Factor: 0.99    Method.....: SW846 8330

PARAMETER	RESULT	REPORTING LIMIT	UNITS	MDL
Nitroglycerin	ND	0.64	ug/L	0.15
PETN	ND	0.64	ug/L	0.23
Picric Acid	ND	0.99	ug/L	0.12
2-Amino-4,6-dinitrotoluene	ND	0.20	ug/L	0.099
4-Amino-2,6-dinitrotoluene	ND	0.099	ug/L	0.022
1,3-Dinitrobenzene	ND	0.099	ug/L	0.050
2,4-Dinitrotoluene	ND	0.099	ug/L	0.050
2,6-Dinitrotoluene	ND	0.099	ug/L	0.050
HMX	ND	0.099	ug/L	0.027
Nitrobenzene	ND	0.099	ug/L	0.050
2-Nitrotoluene	ND	0.40	ug/L	0.071
3-Nitrotoluene	ND	0.40	ug/L	0.061
4-Nitrotoluene	ND	0.50	ug/L	0.071
RDX	2.8 J	0.099	ug/L	0.064
Tetryl	ND	0.099	ug/L	0.050
1,3,5-Trinitrobenzene	ND	0.099	ug/L	0.031
2,4,6-Trinitrotoluene	ND	0.099	ug/L	0.024

SURROGATE	PERCENT RECOVERY	RECOVERY LIMITS
3,4-Dinitrotoluene	99	(79 - 111)

*UK  
9/6/09*

TestAmerica Portland

Client Sample ID: PSF0979-02

HPLC 23L4MWO2AW

Lot-Sample #...: G9G030156-002    Work Order #...: LF3H51AC    Matrix.....: WATER  
 Date Sampled...: 06/29/09    Date Received...: 07/02/09  
 Prep Date.....: 07/06/09    Analysis Date...: 07/09/09  
 Prep Batch #...: 9187206    *0357*  
 Dilution Factor: 1    Method.....: SW846 8330

PARAMETER	RESULT	REPORTING		
		LIMIT	UNITS	MDL
Nitroglycerin	ND	0.65	ug/L	0.15
PETN	ND	0.65	ug/L	0.23
Picric Acid	ND	1.0	ug/L	0.12
2-Amino-4,6-dinitrotoluene	ND	0.20	ug/L	0.10
4-Amino-2,6-dinitrotoluene	ND	0.10	ug/L	0.022
1,3-Dinitrobenzene	ND	0.10	ug/L	0.050
2,4-Dinitrotoluene	ND	0.10	ug/L	0.050
2,6-Dinitrotoluene	ND	0.10	ug/L	0.050
HMX	3.8 J	0.10	ug/L	0.027
Nitrobenzene	ND	0.10	ug/L	0.050
2-Nitrotoluene	ND	0.40	ug/L	0.072
3-Nitrotoluene	ND	0.40	ug/L	0.062
4-Nitrotoluene	ND	0.50	ug/L	0.072
ROX	1.9 J	0.10	ug/L	0.065
Tetryl	ND	0.10	ug/L	0.050
1,3,5-Trinitrobenzene	ND	0.10	ug/L	0.031
2,4,6-Trinitrotoluene	ND	0.10	ug/L	0.024
		PERCENT	RECOVERY	
<u>SURROGATE</u>	<u>RECOVERY</u>	<u>LIMITS</u>		
3,4-Dinitrotoluene	102	(79 - 111)		

*WV  
8/10/09*



TestAmerica Portland

Client Sample ID: PSF0979-03

HPLC 23L4MW02BW

Lot-Sample #...: G9G030156-003 Work Order #...: LF3H61AC Matrix.....: WATER  
 Date Sampled...: 06/29/09 Date Received...: 07/02/09  
 Prep Date.....: 07/06/09 Analysis Date...: 07/11/09  
 Prep Batch #...: 9187206  
 Dilution Factor: 5 Method.....: SW846 8330

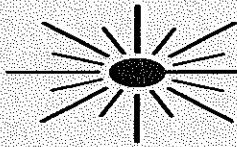
PARAMETER	RESULT	REPORTING		
		LIMIT	UNITS	MDL
Nitroglycerin	ND	3.2	ug/L	0.75
PETN	ND	3.2	ug/L	1.2
Picric Acid	ND	5.0	ug/L	0.62
2-Amino-4,6-dinitrotoluene	ND	1.0	ug/L	0.50
4-Amino-2,6-dinitrotoluene	ND	0.50	ug/L	0.11
1,3-Dinitrobenzene	ND	0.50	ug/L	0.25
2,4-Dinitrotoluene	0.36 J	0.50	ug/L	0.25
2,6-Dinitrotoluene	ND	0.50	ug/L	0.25
HMX	4.3 J	0.50	ug/L	0.14
Nitrobenzene	ND	0.50	ug/L	0.25
2-Nitrotoluene	ND	2.0	ug/L	0.36
3-Nitrotoluene	ND	2.0	ug/L	0.31
4-Nitrotoluene	ND	2.5	ug/L	0.36
RDX	84 J	0.50	ug/L	0.32
Tetryl	ND	0.50	ug/L	0.25
1,3,5-Trinitrobenzene	ND	0.50	ug/L	0.16
2,4,6-Trinitrotoluene	ND	0.50	ug/L	0.12

SURROGATE	PERCENT	RECOVERY
	RECOVERY	LIMITS
3,4-Dinitrotoluene	102	(79 - 111)

NOTE(S):

J Estimated result. Result is less than RL.

WJ  
8/10/09



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## *Support Documentation*

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*Laboratory Case Narrative*

## **DATA DELIVERABLES PACKAGE**

Michael Baker Jr., Inc.  
James D. Peyton  
5261 Fountain Drive, Suite A  
Crown Point, IN 46307

Client Project: Camp Bonneville Groundwater  
Client Project Number: 110002 Task 6200

TA Work Order#: PSF0979  
TA Project Manager: Estella Rieben

The total number of pages contained in this data package is:

147

July 31, 2009

TestAmerica - Portland  
9405 S. W. Nimbus Avenue  
Beaverton, Oregon 97008  
(503) 906-9200  
(503) 906-9210

## CASE NARRATIVE

**Client:** Michael Baker Jr., Inc. **Date Sampled:** 6/29/2009  
**Project:** Camp Bonneville Groundwater **Date Received:** 6/30/2009  
110002 Task 6200  
**Lab:** PSF0979

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**SAMPLE RECEIPT:** Samples were received intact, on ice, with chain of custody documentation. The sample temperature was measured at 2.3° C upon receipt at the laboratory.

**HOLDING TIMES:** All samples were analyzed within holding times.

**PROBLEMS ENCOUNTERED:** No problems were encountered.

**QA/QC CRITERIA:** No problems were encountered during sample analysis.

**OBSERVATIONS:** No significant observations were made.

**SUBCONTRACTED:** 314.0\_9196314 (Perchlorate) and SW846 8330 analyses were performed at the TestAmerica facility in West Sacramento, CA. The data package for the specified analyses is included in this data deliverables package as Appendix A.

*"I certify that this data package is in compliance with the contract both technically and for completeness, for all conditions other than the conditions detailed above. Release of the data contained in this data package has been authorized by the Laboratory Director or his designee, as verified by the following signature."*

TestAmerica Portland



Estella Rieben  
Project Manager

**Michael Baker Jr., Inc.**  
5261 Fountain Drive, Suite A  
Crown Point, IN 46307

Project Name: **Camp Bonneville Groundwater**  
Project Number: 110002 Task 6200  
Project Manager: James D. Peyton

Report Created:  
07/21/09 08:55

## ANALYTICAL REPORT FOR SAMPLES

Sample ID	Laboratory ID	Matrix	Date Sampled	Date Received
23L4MW04AW	PSF0979-01	Water	06/29/09 16:30	06/30/09 17:10
23L4MW02AW	PSF0979-02	Water	06/29/09 18:00	06/30/09 17:10
23L4MW02BW	PSF0979-03	Water	06/29/09 21:00	06/30/09 17:10
TB-261	PSF0979-04	Water	06/29/09 00:00	06/30/09 17:10

TestAmerica Portland



Estella Rieben, Project Manager

*The results in this report apply to the samples analyzed in accordance with the chain of custody document. This analytical report shall not be reproduced except in full, without the written approval of the laboratory.*

Michael Baker Jr., Inc.

5261 Fountain Drive, Suite A  
Crown Point, IN 46307

Project Name: **Camp Bonneville Groundwater**

Project Number: 110002 Task 6200

Project Manager: James D. Peyton

Report Created:

07/21/09 08:55

## Notes and Definitions

### Report Specific Notes:

- a - Spiked analyte recovery is outside stated control limits.
- J - Estimated result. Result is less than RL.
- Q - Elevated Reporting limit. The reporting limit is elevated due to high analyte levels.

### Laboratory Reporting Conventions:

- DET - Analyte DETECTED at or above the Reporting Limit. Qualitative Analyses only.
- ND - Analyte NOT DETECTED at or above the reporting limit (MDL or MRL, as appropriate).
- NR/NA - Not Reported / Not Available
- dry - Sample results reported on a Dry Weight Basis. Results and Reporting Limits have been corrected for Percent Dry Weight.
- wet - Sample results and reporting limits reported on a Wet Weight Basis (as received). Results with neither 'wet' nor 'dry' are reported on a Wet Weight Basis.
- RPD - RELATIVE PERCENT DIFFERENCE (RPDs calculated using Results, not Percent Recoveries).
- MRL - METHOD REPORTING LIMIT. Reporting Level at, or above, the lowest level standard of the Calibration Table.
- MDL\* - METHOD DETECTION LIMIT. Reporting Level at, or above, the statistically derived limit based on 40CFR, Part 136, Appendix B. \*MDLs are listed on the report only if the data has been evaluated below the MRL. Results between the MDL and MRL are reported as Estimated Results.
- Dil - Dilutions are calculated based on deviations from the standard dilution performed for an analysis, and may not represent the dilution found on the analytical raw data.
- Reporting Limits - Reporting limits (MDLs and MRLs) are adjusted based on variations in sample preparation amounts, analytical dilutions and percent solids, where applicable.
- Electronic Signature - Electronic Signature added in accordance with TestAmerica's *Electronic Reporting and Electronic Signatures Policy*. Application of electronic signature indicates that the report has been reviewed and approved for release by the laboratory. Electronic signature is intended to be the legally binding equivalent of a traditionally handwritten signature.

TestAmerica Portland



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## Case Narrative

### TestAmerica West Sacramento Project Number G9G030156

#### **WATER, 8330, Nitroaromatics & Nitramines**

Samples: 1, 2, 3

The laboratory control sample (LCS) associated with this extraction batch has percent recoveries for some analytes above the established control limits indicating a potential high bias in the data. There is no adverse impact upon the data for the analytes that are not detected at concentrations above the reporting limit in the samples. The samples do have detected concentrations of RDX and HMX but this was not discovered until after the recommended hold time for extraction had expired. After consulting with Ms. Rieben, the samples were not re-extracted and the analytical results are included in this report.

There was insufficient sample volume to prepare a matrix spike/matrix spike duplicate (MS/MSD) pair with this extraction batch.

There are no other anomalies associated with this project.



## Sample Summary

### TestAmerica West Sacramento Project Number G9G030156

<u>WO#</u>	<u>Sample #</u>	<u>Client Sample ID</u>	<u>Sampling Date</u>	<u>Received Date</u>
LF3H4	1	PSF0979-01 234MWO4AW	6/29/2009 04:30 PM	7/2/2009 09:50 AM
LF3H5	2	PSF0979-02 234MWO2BW	6/29/2009 06:00 PM	7/2/2009 09:50 AM
LF3H6	3	PSF0979-03 234MWO2BW	6/29/2009 09:00 PM	7/2/2009 09:50 AM

#### Notes(s):

- The analytical results of the samples listed above are presented on the following pages.
- All calculations are performed before rounding to avoid round-off errors in calculated results.
- Results noted as "ND" were not detected at or above the stated limit.
- This report must not be reproduced, except in full, without the written approval of the laboratory.
- Results for the following parameters are never reported on a dry weight basis: color, corrosivity, density, flashpoint, ignitability, layers, odor, paint filter test, pH, porosity, pressure, reactivity, redox potential, specific gravity, spot tests, solids, solubility, temperature, viscosity, and weight.

# QC DATA ASSOCIATION SUMMARY

G9G030156

Sample Preparation and Analysis Control Numbers

<u>SAMPLE#</u>	<u>MATRIX</u>	<u>ANALYTICAL METHOD</u>	<u>LEACH BATCH #</u>	<u>PREP BATCH #</u>	<u>MS RUN#</u>
001	WATER	SW846 8330		9187206	
002	WATER	SW846 8330		9187206	
003	WATER	SW846 8330		9187206	

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## *Chain of Custody*



TestAmerica Portland  
**Sample Receiving Checklist**

Work Order #: PSF0979 Date/Time Received: 6/30/09 1710  
 Client Name and Project: Michael Baker

Time Zone:  
 EDT/EST     CDT/CST     MDT/MST     PDT/PST     AK     OTHER

**Unpacking Checks:**

Cooler #(s): \_\_\_\_\_  
 Temperatures: \_\_\_\_\_  
                   Digi #1 Digi #2 IR Gun  
                                  ( Plastic  Glass)

**Temperature out of Range:**

\_\_\_\_ Not enough or No Ice  
 \_\_\_\_ Ice Melted  
 \_\_\_\_ W/in 4 Hrs of collection  
 \_\_\_\_ Other: \_\_\_\_\_

N/A    Yes    No

Initials: BLE

- 1. If ESI client, were temp blanks received? If no, document on NOD.
- 2. Cooler Seals intact? (N/A if hand delivered) if no, document on NOD.
- 3. Chain of Custody present? If no, document on NOD.
- 4. Bottles received intact? If no, document on NOD.
- 5. Sample is not multiphasic? If no, document on NOD.
- 6. Proper Container and preservatives used? If no, document on NOD.
- 7. pH of all samples checked and meet requirements? If no, document on NOD.
- 8. Cyanide samples checked for sulfides and meet requirements? If no, notify PM.
- 9. HF Dilution required?
- 10. Sufficient volume provided for all analysis? If no, document on NOD and consult PM before proceeding.
- 11. Did chain of custody agree with samples received? If no, document on NOD.
- 12. Is the "Sampled by" section of the COC completed?
- 13. Were VOA/Oil Syringe samples without headspace?
- 14. Were VOA vials preserved?  HCl     Sodium Thiosulfate     Ascorbic Acid
- 15. Did samples require preservation with sodium thiosulfate?
- 16. If yes to #14, was the residual chlorine test negative? If no, document on NOD.
- 17. Are dissolved/field filtered metals bottles sediment-free? If no, document on NOD.
- 18. Is sufficient volume provided for client requested MS/MSD or matrix duplicates? If no, document on NOD and contact PM before proceeding.
- 19. Are analyses with short holding times received in hold?
- 20. Was Standard Turn Around (TAT) requested?
- 21. Receipt date(s) < 48 hours past the collection date(s)? If no, notify PM.

TestAmerica Portland  
Sample Receiving Checklist

Work Order #: PSFO979

**Login Checks:**

Initials: PS

N/A Yes No

22. Sufficient volume provided for all analysis? If no, document on NOD & contact PM.
23. Sufficient volume provided for client requested MS/MSD or matrix duplicates? If no, document on NOD and contact PM.
24. Did the chain of custody include "received by" and "relinquished by" signatures, dates and times?
25. Were special log in instructions read and followed?
26. Were tests logged checked against the COC?
27. Were rush notices printed and delivered?
28. Were short hold notices printed and delivered?
29. Were subcontract COCs printed?
30. Was HF dilution logged?

**Labeling and Storage Checks:**

Initials: PS

N.A Yes No

31. Were the subcontracted samples/containers put in Sx fridge?
32. Were sample bottles and COC double checked for dissolved/filtered metals?
33. Did the sample ID, Date, and Time from label match what was logged?
34. Were Foreign sample stickers affixed to each container and containers stored in foreign fridge?
35. Were HF stickers affixed to each container, and containers stored in Sx fridge?
36. Was an NOD for created for noted discrepancies and placed in folder?

Document any problems or discrepancies and the actions taken to resolve them on a Notice of Discrepancy form (NOD).

CLIENT TAL - Portland PM LL LOG# 99325  
LOT# (QUANTIMS ID) 996030154 QUOTE# 87265 LOCATION W25C

DATE RECEIVED 7-2-09 TIME RECEIVED 950 Initials AK Date 7-2-09

DELIVERED BY  FEDEX  CA OVERNIGHT  CLIENT  
 AIRBORNE  GOLDENSTATE  DHL  
 UPS  BAX GLOBAL  GO-GETTERS  
 TAL COURIER  VALLEY LOGISTICS  MORGAN HILL COURIER  
 OTHER

CUSTODY SEAL STATUS  INTACT  BROKEN  N/A

CUSTODY SEAL #(S) 204635, 204625

SHIPPING CONTAINER(S)  TAL  CLIENT  N/A

TEMPERATURE RECORD (IN °C) IR 4  5  OTHER

COC #(S) AK

TEMPERATURE BLANK Observed: 0 Corrected: 2

SAMPLE TEMPERATURE Observed: 1 2 2 Average: 2 Corrected Average: 2

COLLECTOR'S NAME:  Verified from COC  Not on COC

pH MEASURED  YES  ANOMALY  N/A

LABELED BY .....

LABELS CHECKED BY .....

PEER REVIEW  NA

SHORT HOLD TEST NOTIFICATION SAMPLE RECEIVING  
WETCHEM  N/A  
VOA-ENCORES  N/A

METALS NOTIFIED OF FILTER/PRESERVE VIA VERBAL & EMAIL  N/A

COMPLETE SHIPMENT RECEIVED IN GOOD CONDITION WITH APPROPRIATE TEMPERATURES, CONTAINERS, PRESERVATIVES  N/A

CLOUSEAU  TEMPERATURE EXCEEDED (2 °C - 6 °C)<sup>\*1</sup>  N/A

WET ICE  BLUE ICE  GEL PACK  NO COOLING AGENTS USED  PM NOTIFIED

Notes: \_\_\_\_\_

\*1 Acceptable temperature range for State of Wisconsin samples is ≤4°C

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# *Worksheets*



**HOLDING TIMES**

SAMPLE DATE	SAMPLE ID	Portland LAB ID	Sacramento Lab ID	Matrix	Prep	Anal
06/29/09	23L4MW04AW	PSF0979-01	LF3H4	Aq	07/06/09	7/9 1X
06/29/09	23L4MW02AW	PSF0979-02	LF3H5	Aq	07/06/09	7/9 1X
06/29/09	23L4MW02BW	PSF0979-03	LF3H6	Aq	07/06/09	7/11 5X

7 days from collection to extraction  
 Sample Date 6/29/2009 Extract By 7/6/2009  
 40 days from extraction to analysis  
 Sample Date 7/6/2009 Extract By 7/13/2009

**TARGET COMPOUNDS AND QUANTITATION LIMITS**

Was a target compound list provided by the client? No  
 Did Sample Form 1s match the target compound list NA  
 Were required quantitation limits provided by the client? No  
 Did all compounds meet the required quantitation limits? NA

**SYSTEM MONITORING COMPOUNDS**

SAMPLE	SURROGATE	COLUMN 1	DF	ACTION
All IN				

Were surrogate RTs within windows established by the ICAL? Yes  
 Were there any transcription errors between the raw data and Form 2? No  
 Were laboratory acceptance limits used as the basis for validation? 30-150  
 Did the laboratory provide CLP Form II or equivalent? Yes

02BW

SURR	AMOUNT FOUND	AMOUNT SPIKED	% R	F1
34-DNT A	2.535	2.5	101.4	102
34-DNT C	2.952	2.5	118.1	

**MS/MSD RECOVERY and RELATIVE PERCENT DIFFERENCE**

None Present

**LABORATORY CONTROL SAMPLES**

SAMPLE	COMPOUND	LCS %R	LCSD %R	RPD	ACTION
LF3161A	4-a-26-DNTol	114	NA	NA	68-113
	HMX	116	NA	NA	67-115
	RDX	124	NA	NA	68-122
	135-TNTol	136	NA	NA	74-120
			NA	NA	

Were laboratory acceptance limits used as the basis for validation? Yes  
 Did the laboratory provide CLP Form III or equivalent? Yes  
 Were chromatograms and quan reports present for all LCS/LCDs? Yes

LF3161A

4-a-26-DNTol	AMOUNT FOUND	AMOUNT SPIKED	% R	FORM 3
LCS	1.14	1	114.0	114

%R = (Amount Found/Amount Spiked)\*100

**BLANKS**

BLANK	COMPOUND	RESULT	5X OR 10X	ACTION LEVEL	ACTION
LF3161A	All Parameters	ND		#VALUE!	

NOTE Equipment/Field blanks are not qualified on the basis of laboratory method blank contamination or contamination in other field quality control blanks.

**CALIBRATION**

**Initial**

Are chromatograms and quan reports present for all ICAL standards? Yes  
 Are CLP For VIs or equivalent present and complete? Yes  
 Any transcription or calculation errors? No  
 What linearity criteria was used? 20% RSD or 0.99  
 Were RT window documented? Yes

**Continuing**

Are chromatograms and quan reports present for all CCV standards? Yes  
 Are CLP Form VIIIs or equivalent present and complete? Yes  
 Any transcription or calculation errors? No  
 What %D criteria was used? 25% D  
 Have all samples and standards been listed on an Analytical Sequence? Yes  
 Was a proper analytical sequence followed? Yes

**INITIAL CALIBRATION - A**

Date	6/10/2009	Rptd Avg CF	77.89163	Rptd %RSD	5.156
Instrument	A	Calcd Avg CF	77.892	Calcd %RSD	5.156
Compound	RDX				
Reported CF	80.13	CF1	81.80	STD DEV	4
Calculated CF	80.42	CF2	75.90		
		CF3	77.15		
Response	16084	CF4	80.72		
Conc	200	CF5	78.45		
		CF6	80.13		
	CF = (H/C)	CF7	69.168	%RSD = (Std Dev/Avg CF)*100	
		CF8	79.815		

**CONTINUING CALIBRATION A**

Date	7/9/2009				
Time	0035				
Instrument	A				
Compound	HMX				
Reported Conc	104.5	Rptd %D	5		
Calculated Conc	104.380	Calcd %D	-4.38		
Response	12630	ICAL CF	100		
CF	121	CCV CF	104.4		
	CF = (H/C)			%Difference = ((ICAL - CCV)/ICAL)*100	

**INITIAL CALIBRATION - C**

Date	6/9/2009	Rptd Avg CF	42.55812	Rptd %RSD	4.852
Instrument	C	Calcd Avg CF	42.558	Calcd %RSD	4.852
Compound	RDX				
Reported CF	42.585	CF1	45.40	STD DEV	2
Calculated CF	42.78	CF2	43.70		
		CF3	42.75		
Response	8556	CF4	43.30		
Conc	200	CF5	42.16		
		CF6	42.59		
	CF = (H/C)	CF7	38.118	%RSD = (Std Dev/Avg CF)*100	
		CF8	42.452		

**CONTINUING CALIBRATION C**

Date	7/11/2009				
Time	0554				
Instrument	C				
Compound	HMX				
Reported Conc	184.3	Rptd %D	8		
Calculated Conc	184.280	Calcd %D	7.86		
Response	7594	ICAL CF	200		
CF	41.20913	CCV CF	184.3		
	CF = (H/C)			%Difference = ((ICAL - CCV)/ICAL)*100	

**FIELD DUPLICATES**

COMPOUND	SAMPLE	QUALIFIER	DUPLICATE	QUALIFIER	RPD
PCA					#DIV/0!

NOTES      Samples are not qualified on this basis.

**COMPOUND IDENTIFICATION AND QUANTITATION**

Has a F10 been completed for every sample containing positive results?	Yes
Was RT data presented on the form?	NA
Are RTs within the established windows?	Yes
Any transcription or calculation errors?	No
Any false positives, negative peaks, shouldering, etc.?	No
Was GCMS confirmation needed for results > 10 µg/ml?	NA
Were percent differences or relative percent differences calculated?	NA
Are percent differences/RPDs greater than 25%?	NA
Are there any transcription errors?	No
Are Form 1s present for all field and quality control samples?	No
Are chromatograms and quan reports present for all samples?	Yes
Are RLs adjusted to reflect sample dilutions, percent solids, etc.?	Yes
For soils, any percent solids <50%?	No
For soils, any percent solids <10%?	No

23L4MW02BW RDX 84 µg/L

	A	C
response	65192	31840
cf	77.89163	42.55812
final vol ml	20	20
initial L	1.0053	1.0053
df	5	5
calculated	83.2545	74.4209

**PERCENT DIFFERENCE BETWEEN COLUMNS**

SAMPLE	COMPOUND	A	C	RPD	QUALIFIER
23L4MW04AW	RDX	2.814	2.589	8.3	none
23L4MW02AW	HMX	3.825	3.335	13.7	none
	RDX	19.19	17.52	9.1	none
23L4MW02BW	24-DNT	0.3578	0.4359	-19.7	none
	HMX	4.265	3.572	17.7	none
	RDX	83.25	74.42	11.2	none

<b>Michael Baker Jr., Inc.</b> 5261 Fountain Drive, Suite A Crown Point, IN 46307	<b>Project Name:</b> Camp Bonneville Groundwater <b>Project Number:</b> I10002 Task 6200 <b>Project Manager:</b> James D. Peyton	<b>Report Created:</b> 07/21/09 08:55
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**SW846 8330**  
TestAmerica West Sacramento

Analyte	Method	Result	MDL*	MRL	Units	Dil	Batch	Prepared	Analyzed	Notes
<b>PSF0979-01 (23L4MW04AW)</b>		<b>Water</b>				<b>Sampled: 06/29/09 16:30</b>				
1,3,5-Trinitrobenzene	SW846 8330	ND	0.031	0.099	ug/L	0.99x	9187206	07/06/09 09:30	07/09/09 03:06	
1,3-Dinitrobenzene	"	ND	0.05	0.099	"	"	"	"	"	
2,4,6-Trinitrotoluene	"	ND	0.024	0.099	"	"	"	"	"	
2,4-Dinitrotoluene	"	ND	0.05	0.099	"	"	"	"	"	
2,6-Dinitrotoluene	"	ND	0.05	0.099	"	"	"	"	"	
2-Amino-4,6-dinitrotoluene	"	ND	0.099	0.2	"	"	"	"	"	
2-Nitrotoluene	"	ND	0.071	0.4	"	"	"	"	"	
3-Nitrotoluene	"	ND	0.061	0.4	"	"	"	"	"	
4-Amino-2,6-dinitrotoluene	"	ND	0.022	0.099	"	"	"	"	"	
4-Nitrotoluene	"	ND	0.071	0.5	"	"	"	"	"	
HMX	"	ND	0.027	0.099	"	"	"	"	"	
Nitrobenzene	"	ND	0.05	0.099	"	"	"	"	"	
Nitroglycerin	"	ND	0.15	0.64	"	"	"	"	"	
PETN	"	ND	0.23	0.64	"	"	"	"	"	
Picric Acid	"	ND	0.12	0.99	"	"	"	"	"	
<b>RDX</b>	"	<b>2.8</b>	<b>0.064</b>	<b>0.099</b>	"	"	"	"	"	
Tetryl	"	ND	0.05	0.099	"	"	"	"	"	
<i>Surrogate(s): 3,4-Dinitrotoluene</i>			99%	79 - 111 %	"	"	"	"	"	"

<b>PSF0979-02 (23L4MW02AW)</b>		<b>Water</b>				<b>Sampled: 06/29/09 18:00</b>				
1,3,5-Trinitrobenzene	SW846 8330	ND	0.031	0.1	ug/L	1x	9187206	07/06/09 09:30	07/09/09 03:57	
1,3-Dinitrobenzene	"	ND	0.05	0.1	"	"	"	"	"	
2,4,6-Trinitrotoluene	"	ND	0.024	0.1	"	"	"	"	"	
2,4-Dinitrotoluene	"	ND	0.05	0.1	"	"	"	"	"	
2,6-Dinitrotoluene	"	ND	0.05	0.1	"	"	"	"	"	
2-Amino-4,6-dinitrotoluene	"	ND	0.1	0.2	"	"	"	"	"	
2-Nitrotoluene	"	ND	0.072	0.4	"	"	"	"	"	
3-Nitrotoluene	"	ND	0.062	0.4	"	"	"	"	"	
4-Amino-2,6-dinitrotoluene	"	ND	0.022	0.1	"	"	"	"	"	
4-Nitrotoluene	"	ND	0.072	0.5	"	"	"	"	"	
<b>HMX</b>	"	<b>3.8</b>	<b>0.027</b>	<b>0.1</b>	"	"	"	"	"	
Nitrobenzene	"	ND	0.05	0.1	"	"	"	"	"	
Nitroglycerin	"	ND	0.15	0.65	"	"	"	"	"	
PETN	"	ND	0.23	0.65	"	"	"	"	"	
Picric Acid	"	ND	0.12	1	"	"	"	"	"	
<b>RDX</b>	"	<b>19</b>	<b>0.065</b>	<b>0.1</b>	"	"	"	"	"	

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*Estella K Rieben*

Estella Rieben, Project Manager

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**SW846 8330**  
TestAmerica West Sacramento

Analyte	Method	Result	MDL*	MRL	Units	Dil	Batch	Prepared	Analyzed	Notes
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**PSF0979-02 (23L4MW02AW)** Water Sampled: 06/29/09 18:00

Tetryl	SW846 8330	ND	0.05	0.1	ug/L	1x	9187206	07/06/09 09:30	07/09/09 03:57	
<i>Surrogate(s): 3,4-Dinitrotoluene</i>			102%			79 - 111 %	"			"

**PSF0979-03 (23L4MW02BW)** Water Sampled: 06/29/09 21:00

1,3,5-Trinitrobenzene	SW846 8330	ND	0.16	0.5	ug/L	5x	9187206	07/06/09 09:30	07/11/09 06:11	
1,3-Dinitrobenzene	"	ND	0.25	0.5	"	"	"	"	"	
2,4,6-Trinitrotoluene	"	ND	0.12	0.5	"	"	"	"	"	
<b>2,4-Dinitrotoluene</b>	"	<b>0.36</b>	0.25	0.5	"	"	"	"	"	<b>J</b>
2,6-Dinitrotoluene	"	ND	0.25	0.5	"	"	"	"	"	
2-Amino-4,6-dinitrotoluene	"	ND	0.5	1	"	"	"	"	"	
2-Nitrotoluene	"	ND	0.36	2	"	"	"	"	"	
3-Nitrotoluene	"	ND	0.31	2	"	"	"	"	"	
4-Amino-2,6-dinitrotoluene	"	ND	0.11	0.5	"	"	"	"	"	
4-Nitrotoluene	"	ND	0.36	2.5	"	"	"	"	"	
<b>HMX</b>	"	<b>4.3</b>	0.14	0.5	"	"	"	"	"	
Nitrobenzene	"	ND	0.25	0.5	"	"	"	"	"	
Nitroglycerin	"	ND	0.75	3.2	"	"	"	"	"	
PETN	"	ND	1.2	3.2	"	"	"	"	"	
Picric Acid	"	ND	0.62	5	"	"	"	"	"	
<b>RDX</b>	"	<b>84</b>	0.32	0.5	"	"	"	"	"	
Tetryl	"	ND	0.25	0.5	"	"	"	"	"	
<i>Surrogate(s): 3,4-Dinitrotoluene</i>			102%			79 - 111 %	"			"

TestAmerica Portland

*Estella K. Rieben*

Estella Rieben, Project Manager

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**SW846 8330 - Laboratory Quality Control Results**  
TestAmerica West Sacramento

QC Batch: 9187206	WATER Preparation Method: NA
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Analyte	Method	Result	MDL*	MRI	Units	Dil	Source Result	Spike Amt	% REC	(Limits)	% RPD	(Limits)	Analyzed	Notes
<b>Blank (G9G060000206B)</b>			QC Source:				Extracted: 07/06/09 09:30							
1,3,5-Trinitrobenzene	SW846 8330	ND	0.031	0.1	ug/L	1x	--	--	--	--	--	--	07/09/09 01:25	
1,3-Dinitrobenzene	"	ND	0.05	0.1	"	"	--	--	--	--	--	--	"	
2,4,6-Trinitrotoluene	"	ND	0.024	0.1	"	"	--	--	--	--	--	--	"	
2,4-Dinitrotoluene	"	ND	0.05	0.1	"	"	--	--	--	--	--	--	"	
2,6-Dinitrotoluene	"	ND	0.05	0.1	"	"	--	--	--	--	--	--	"	
2-Amino-4,6-dinitrotoluene	"	ND	0.1	0.2	"	"	--	--	--	--	--	--	"	
2-Nitrotoluene	"	ND	0.072	0.4	"	"	--	--	--	--	--	--	"	
3-Nitrotoluene	"	ND	0.062	0.4	"	"	--	--	--	--	--	--	"	
4-Amino-2,6-dinitrotoluene	"	ND	0.022	0.1	"	"	--	--	--	--	--	--	"	
4-Nitrotoluene	"	ND	0.072	0.5	"	"	--	--	--	--	--	--	"	
HMX	"	ND	0.027	0.1	"	"	--	--	--	--	--	--	"	
Nitrobenzene	"	ND	0.05	0.1	"	"	--	--	--	--	--	--	"	
Nitroglycerin	"	ND	0.15	0.65	"	"	--	--	--	--	--	--	"	
PETN	"	ND	0.23	0.65	"	"	--	--	--	--	--	--	"	
Picric Acid	"	ND	0.12	1	"	"	--	--	--	--	--	--	"	
RDX	"	ND	0.065	0.1	"	"	--	--	--	--	--	--	"	
Tetryl	"	ND	0.05	0.1	"	"	--	--	--	--	--	--	"	
<i>Surrogate(s): 3,4-Dinitrotoluene</i>		<i>Recovery: 102%</i>		<i>Limits: 79-111%</i>				<i>"</i>						<i>07/09/09 01:25</i>

<b>LCS (G9G060000206C)</b>			QC Source:				Extracted: 07/06/09 09:30							
1,3,5-Trinitrobenzene	SW846 8330	1.36	0.031	0.1	ug/L	1x	--	1	136%	(74-120)	--	--	07/09/09 02:16	a
1,3-Dinitrobenzene	"	1.23	0.05	0.1	"	"	--	"	123%	(72-123)	--	--	"	
2,4,6-Trinitrotoluene	"	1.11	0.024	0.1	"	"	--	"	111%	(69-111)	--	--	"	
2,4-Dinitrotoluene	"	1.18	0.05	0.1	"	"	--	"	118%	(70-119)	--	--	"	
2,6-Dinitrotoluene	"	1.19	0.05	0.1	"	"	--	"	119%	(71-119)	--	--	"	
2-Amino-4,6-dinitrotoluene	"	1.22	0.1	0.2	"	"	--	"	122%	(77-123)	--	--	"	
2-Nitrotoluene	"	1.09	0.072	0.5	"	"	--	"	109%	(64-120)	--	--	"	
3-Nitrotoluene	"	1.11	0.062	0.5	"	"	--	"	111%	(67-114)	--	--	"	
4-Amino-2,6-dinitrotoluene	"	1.14	0.022	0.1	"	"	--	"	114%	(68-113)	--	--	"	a
4-Nitrotoluene	"	1.12	0.072	0.5	"	"	--	"	112%	(67-115)	--	--	"	
HMX	"	1.16	0.027	0.1	"	"	--	"	116%	"	--	--	"	a
Nitrobenzene	"	1.19	0.05	0.1	"	"	--	"	119%	(69-119)	--	--	"	
Nitroglycerin	"	5.1	0.15	0.65	"	"	--	5	102%	(85-115)	--	--	"	
PETN	"	5.24	0.23	0.65	"	"	--	"	105%	(84-117)	--	--	"	
Picric Acid	"	3.29	0.12	1	"	"	--	"	66%	(21-118)	--	--	"	
RDX	"	1.24	0.065	0.1	"	"	--	1	124%	(68-122)	--	--	"	a
Tetryl	"	0.986	0.05	0.1	"	"	--	"	99%	(66-105)	--	--	"	
<i>Surrogate(s): 3,4-Dinitrotoluene</i>		<i>Recovery: 101%</i>		<i>Limits: 79-111%</i>				<i>"</i>						<i>07/09/09 02:16</i>

TestAmerica Portland

*Estella K Rieben*

Estella Rieben, Project Manager

*The results in this report apply to the samples analyzed in accordance with the chain of custody document. This analytical report shall not be reproduced except in full, without the written approval of the laboratory.*

LABORATORY CONTROL SAMPLE EVALUATION REPORT

HPLC

Client Lot #...: G9G030156      Work Order #...: LF3161AC      Matrix.....: WATER  
 LCS Lot-Sample#: G9G060000-206  
 Prep Date.....: 07/06/09      Analysis Date...: 07/09/09  
 Prep Batch #...: 9187206  
 Dilution Factor: 1

<u>PARAMETER</u>	<u>PERCENT RECOVERY</u>	<u>RECOVERY LIMITS</u>	<u>METHOD</u>
Nitroglycerin	102	(85 - 115)	SW846 8330
PETN	105	(84 - 117)	SW846 8330
Picric Acid	66	(21 - 118)	SW846 8330
2-Amino-4,6-dinitrotoluene	122	(77 - 123)	SW846 8330
4-Amino-2,6-dinitrotoluene	114 a	(68 - 113)	SW846 8330
1,3-Dinitrobenzene	123	(72 - 123)	SW846 8330
2,4-Dinitrotoluene	113	(70 - 119)	SW846 8330
2,6-Dinitrotoluene	119	(71 - 119)	SW846 8330
HMX	116 a	(67 - 115)	SW846 8330
Nitrobenzene	119	(69 - 119)	SW846 8330
2-Nitrotoluene	109	(64 - 120)	SW846 8330
3-Nitrotoluene	111	(67 - 114)	SW846 8330
4-Nitrotoluene	112	(67 - 115)	SW846 8330
RDX	124 a	(68 - 122)	SW846 8330
Tetryl	99	(66 - 105)	SW846 8330
1,3,5-Trinitrobenzene	136 a	(74 - 120)	SW846 8330
2,4,6-Trinitrotoluene	111	(69 - 111)	SW846 8330

<u>SURROGATE</u>	<u>PERCENT RECOVERY</u>	<u>RECOVERY LIMITS</u>
3,4-Dinitrotoluene	101	(79 - 111)

NOTE(S):

Calculations are performed before rounding to avoid round-off errors in calculated results.

Bold print denotes control parameters

a Spiked analyte recovery is outside stated control limits.

LABORATORY CONTROL SAMPLE DATA REPORT

HPLC

Client Lot #...: G9G030156      Work Order #...: LF3161AC      Matrix.....: WATER  
 LCS Lot-Sample#: G9G060000-206  
 Prep Date.....: 07/06/09      Analysis Date...: 07/09/09  
 Prep Batch #...: 9187206  
 Dilution Factor: 1

<u>PARAMETER</u>	<u>SPIKE</u> <u>AMOUNT</u>	<u>MEASURED</u> <u>AMOUNT</u>	<u>UNITS</u>	<u>PERCENT</u> <u>RECOVERY</u>	<u>METHOD</u>
Nitroglycerin	5.00	5.10	ug/L	102	SW846 8330
PEFN	5.00	5.24	ug/L	105	SW846 8330
Picric Acid	5.00	3.29	ug/L	66	SW846 8330
2-Amino-4,6-dinitrotoluene	1.00	1.22	ug/L	122	SW846 8330
4-Amino-2,6-dinitrotoluene	1.00	1.14 a	ug/L	114	SW846 8330
1,3-Dinitrobenzene	1.00	1.23	ug/L	123	SW846 8330
2,4-Dinitrotoluene	1.00	1.18	ug/L	118	SW846 8330
2,6-Dinitrotoluene	1.00	1.19	ug/L	119	SW846 8330
HMX	1.00	1.16 a	ug/L	116	SW846 8330
Nitrobenzene	1.00	1.19	ug/L	119	SW846 8330
2-Nitrotoluene	1.00	1.09	ug/L	109	SW846 8330
3-Nitrotoluene	1.00	1.11	ug/L	111	SW846 8330
4-Nitrotoluene	1.00	1.12	ug/L	112	SW846 8330
RDX	1.00	1.24 a	ug/L	124	SW846 8330
Tetryl	1.00	0.986	ug/L	99	SW846 8330
1,3,5-Trinitrobenzene	1.00	1.36 a	ug/L	136	SW846 8330
2,4,6-Trinitrotoluene	1.00	1.11	ug/L	111	SW846 8330

<u>SURROGATE</u>	<u>PERCENT</u> <u>RECOVERY</u>	<u>RECOVERY</u> <u>LIMITS</u>
3,4-Dinitrotoluene	101	(79 - 111)

**NOTE (S) :**

Calculations are performed before rounding to avoid round-off errors in calculated results.

Bold print denotes control parameters

a Spiked analyte recovery is outside stated control limits.



METHOD BLANK REPORT

HPLC

Client Lot #...: G9G030156      Work Order #...: LF3161AA      Matrix.....: WATER  
 MB Lot-Sample #: G9G060000-206  
 Analysis Date...: 07/09/09      Prep Date.....: 07/06/09  
 Dilution Factor: 1              Prep Batch #...: 9187206

PARAMETER	RESULT	REPORTING		
		LIMIT	UNITS	METHOD
Nitroglycerin	ND	0.65	ug/L	SW846 8330
PETN	ND	0.65	ug/L	SW846 8330
Picric Acid	ND	1.0	ug/L	SW846 8330
2-Amino-4,6-dinitrotoluene	ND	0.20	ug/L	SW846 8330
4-Amino-2,6-dinitrotoluene	ND	0.10	ug/L	SW846 8330
1,3-Dinitrobenzene	ND	0.10	ug/L	SW846 8330
2,4-Dinitrotoluene	ND	0.10	ug/L	SW846 8330
2,6-Dinitrotoluene	ND	0.10	ug/L	SW846 8330
HMX	ND	0.10	ug/L	SW846 8330
Nitrobenzene	ND	0.10	ug/L	SW846 8330
2-Nitrotoluene	ND	0.40	ug/L	SW846 8330
3-Nitrotoluene	ND	0.40	ug/L	SW846 8330
4-Nitrotoluene	ND	0.50	ug/L	SW846 8330
RDX	ND	0.10	ug/L	SW846 8330
Tetryl	ND	0.10	ug/L	SW846 8330
1,3,5-Trinitrobenzene	ND	0.10	ug/L	SW846 8330
2,4,6-Trinitrotoluene	ND	0.10	ug/L	SW846 8330

SURROGATE	PERCENT	RECOVERY
	RECOVERY	LIMITS
3,4-Dinitrotoluene	102	(79 - 111)

NOTE(S):

Calculations are performed before rounding to avoid round-off errors in calculated results.



RQC058

TestAmerica Laboratories, Inc.  
EXTRACTION BENCH WORKSHEET

Run Date: 7/08/09  
Time: 11:09:59

LEV 1  
EXP 1  
Y 1  
Y 2  
MS/MSD 2

Blank  
Check  
Weights/Volumes  
Spike & Surrogate Worksheet  
Vial contains correct volume  
Labels, greenbars, worksheets  
computer batch: correct & all match  
Anomalies to Extraction Method

Expanded Deliverable  
COC Completed  
Bench Sheet Copied  
Package Submitted to Analytical Group  
Bench Sheet Copied per COC

Extractionist: 002448 Tuan Q. Phan

Concentrationist: 002448 Tuan Q. Phan

\*\*\*\*\*  
\* QC BATCH: 9187206 \*  
\* PREP DATE: 7/06/09 9:30  
\* COMP DATE: 7/08/09 10:00  
\*\*\*\*\*

Reviewer/Date: PHANT / 7/08/09

Nitroaromatics & Nitramines: Explosives (8330)  
SOLID PHASE EXTRACTION (NOMINAL)

EXTR	ANL DUE	LOT# MSRUN# / WORK ORDER	TEST FLGS	EXT MTR	MATRIX	INIT/FLN WT/VOL	PHAS INIT ADJ1	ADJ2	EXTRACTION VOL	EXCHANGE VOL	SOLVENTS	SPIKE STANDARD/ SURROGATE ID	
7/06/09	7/15/09	LF3B4-1-AC	R	B7	A0	WATER	1007.3mL 20.00mL	NA	NA	HOAN/ACN	4.5 HOAC/ACN	5.0	50UL-09GCSV172
COMMENTS: G9G030156-001													
7/06/09	7/15/09	LF3B5-1-AC	R	B7	A0	WATER	996.15mL 20.00mL	NA	NA	HOAN/ACN	4.5 HOAC/ACN	5.0	50UL-09GCSV172
COMMENTS: G9G030156-002													
7/06/09	7/15/09	LF3B6-1-AC	R	B7	A0	WATER	1005.3mL 20.00mL	NA	NA	HOAN/ACN	4.5 HOAC/ACN	5.0	50UL-09GCSV172
COMMENTS: G9G030156-003													
7/06/09	0/00/00	LF318-1-AAB		B7	A0	WATER	1000mL 20.00mL	NA	NA	HOAN/ACN	4.5 HOAC/ACN	5.0	50UL-09GCSV172
COMMENTS: G9G060000-206													
7/06/09	0/00/00	LF316-1-ACC		B7	A0	WATER	1000mL 20.00mL	NA	NA	HOAN/ACN	4.5 HOAC/ACN	5.0	50UL-09GCSV172
COMMENTS: G9G060000-206													

LCS SPIKE 20UL-09GCSV0037 100UL-09GCSV0149 100UL-09GCSV0039; 1% HOAC/ACN 2991  
-95; HOAC/H2O 2991-87B; SPE COLUMN WATER LOT 003339096A; SODIUM CHLORIDE LOT  
R070582 MALLINCKRODT; .45 MILLIPORE WATER LOT R5PN29184

R = RUSH  
E = EPA 600  
M = CLIENT REQ MS/MSD  
C = CLP  
D = EXP.DEL.

NUMBER OF WORK ORDERS IN BATCH: 5

TestAmerica West Sacramento  
GC/LC INSTRUMENT LOG

Inst ID: LC10                      Batch ID: 06102009  
Method ; Method 8330              Test : SOP SAC-LC-0009  
ICAL Date: See Calibration Report

Date	Time	Operator	Sample ID	File ID	Vol or Wt	Extract Vol	Diln	Comments
10-JUN-2009	14:42	fhk	Blank	A-000001.	0 g	0 mL	1	
10-JUN-2009	15:33	fhk	STD_1 09GCSV0048 5.0/0/0/ong/m	A-000002.	0 g	0 mL	1	
10-JUN-2009	16:23	fhk	STD_2 09GCSV0049 10/20/10/10ng	A-000003.	0 g	0 mL	1	
10-JUN-2009	17:14	fhk	STD_3 09GCSV0050 20/50/20/20ng	A-000004.	0 g	0 mL	1	Bad Std
10-JUN-2009	18:05	fhk	STD_4 09GCSV0051 50/100/50/50ng	A-000005.	0 g	0 mL	1	
10-JUN-2009	18:56	fhk	STD_5 09GCSV0053 100/200/100/1	A-000006.	0 g	0 mL	1	
10-JUN-2009	19:46	fhk	STD_6 09GCSV0054 200/500/200/2	A-000007.	0 g	0 mL	1	
10-JUN-2009	20:37	fhk	STD_7 09GCSV0055 500/1000/500/	A-000008.	0 g	0 mL	1	
10-JUN-2009	21:28	fhk	STD_8 09GCSV0056 1000/2000/100	A-000009.	0 g	0 mL	1	
10-JUN-2009	22:18	fhk	Blank	A-000010.	0 g	0 mL	1	
10-JUN-2009	23:09	fhk	ICV_6 09GCSV0397 100/200/100/1	A-000011.	0 g	0 mL	1	
11-JUN-2009	00:00	fhk	STD_5 09GCSV0053 100/200/100/1	A-000012.	0 g	0 mL	1	
11-JUN-2009	00:50	fhk	Surrogate 100ng/mL	A-000013.	0 g	0 mL	1	
11-JUN-2009	12:56	fhk	Primer	A-000014.	0 g	0 mL	1	
11-JUN-2009	13:47	fhk	STD_3 09GCSV0050 20/50/20/20ng	A-000015.	0 g	0 mL	1	reinject flow std

Reported

*John Gillog*

TestAmerica West Sacramento

INITIAL CALIBRATION DATA

Start Cal Date : 10-JUN-2009 15:33  
 End Cal Date : 11-JUN-2009 13:47  
 Quant Method : ESTD  
 Origin : Disabled  
 Target Version : 4.14  
 Integrator : Falcon  
 Method file : \\Terastation\share\GCdata\LC10.I\06102009.B\8330AB.M  
 Last Edit : 11-Jun-2009 15:06 kenneyf  
 Curve Type : Average

Calibration File Names:

- Level 1: \\Terastation\share\GCdata\GCdata\LC10.I\06102009.B\A-000002.d
- Level 2: \\Terastation\share\GCdata\GCdata\LC10.I\06102009.B\A-000003.d
- Level 3: \\Terastation\share\GCdata\GCdata\LC10.I\06102009.B\A-000015.d
- Level 4: \\Terastation\share\GCdata\GCdata\LC10.I\06102009.B\A-000005.d
- Level 5: \\Terastation\share\GCdata\GCdata\LC10.I\06102009.B\A-000006.d
- Level 6: \\Terastation\share\GCdata\GCdata\LC10.I\06102009.B\A-000007.d
- Level 7: \\Terastation\share\GCdata\GCdata\LC10.I\06102009.B\A-000008.d
- Level 8: \\Terastation\share\GCdata\GCdata\LC10.I\06102009.B\A-000009.d

Compound	5.000	10.000	20.000	50.000	100.000	200.000	RRP	% RSD
	Level 1	Level 2	Level 3	Level 4	Level 5	Level 6		
	500.000	1000.000						
	Level 7	Level 8						
2 HMX	119 ✓ 116	120 ✓ 123	115 ✓	125 ✓	123 ✓	124 ✓	121	3.066
3 RDX	81.80000 69.16800	75.90000 ✓ 79.81500	77.15000 ✓	80.72000 ✓	78.45000 ✓	80.13000 ✓	77.89163	5.156
4 EGDN	++++ ++++	++++ ++++	++++	++++	++++	++++	++++	++++
5 Picric ACID	++++ 74.53400	++++ 74.87800	91.72000 ✓	85.62000 ✓	84.85500 ✓	84.88200 ✓	82.77483	8.120
6 1,3,5-Trinitrobenzene	161 136	146 147	145	150	147	148	148	4.610
7 1,3-Dinitrobenzene	145 128	142 141	138	142	139	142	140	3.715
8 3,5-Dinitroaniline	97.20000 82.20800	90.80000 89.25000	89.35000	91.80000	88.83000	90.35000	89.97350	4.581
9 TETRYL	97.60000 ✓ 92.43000	95.10000 ✓ 94.40900	77.40000 ✓	96.02000 ✓	94.31000 ✓	95.06000 ✓	92.79113	6.888

*Ask*  
6/11

TestAmerica West Sacramento

INITIAL CALIBRATION DATA

Start Cal Date : 10-JUN-2009 15:33  
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 Quant Method : ESTD  
 Origin : Disabled  
 Target Version : 4.14  
 Integrator : Falcon  
 Method file : \\Terastation\share\GCdata\LC10.T\06102009.B\8330AB.M  
 Last Edit : 11-Jun-2009 15:06 kenneyf  
 Curve Type : Average

Compound	5.000	10.000	20.000	50.000	100.000	200.000	RRF	% RSD
	Level 1	Level 2	Level 3	Level 4	Level 5	Level 6		
	500.000	1000.000						
	Level 7	Level 8						
10 Nitrobenzene	63.80000	62.30000	56.55000	61.44000	61.06000	62.26000		
	56.02600	61.95700					60.67412	4.659
11 Nitroglycerin	+++++	+++++	+++++	+++++	+++++	+++++	+++++	+++++
12 2,4,6-Trinitrotoluene	94.80000	90.20000	81.90000	89.72000	87.04000	87.55500		
	84.22200	87.17800					87.70937	4.401
13 4-AM-2,6-DNT	67.00000	63.30000	60.85000	63.26000	61.61000	62.67000		
	58.65400	61.14600					62.31125	3.901
14 2-AM-4,6-DNT	74.80000	75.80000	73.10000	75.54000	73.53000	74.39000		
	69.18400	72.13500					73.55987	2.932
15 2,6-Dinitrotoluene	57.60000	53.30000	51.90000	53.68000	53.03000	53.21500		
	50.43000	52.74900					53.23800	3.835
16 2,4-Dinitrotoluene	93.00000	86.50000	82.60000	86.20000	84.72000	85.49000		
	81.43200	85.05700					85.62488	4.031
17 2-Nitrotoluene	40.80000	37.40000	33.20000	35.64000	35.21000	35.77500		
	33.97400	35.45000					35.93113	6.486
18 4-Nitrotoluene	48.00000	46.40000	40.75000	43.56000	43.22000	43.56500		
	41.67600	43.21300					43.79800	5.390
19 3-Nitrotoluene	47.00000	44.60000	39.85000	43.24000	42.93000	43.08000		
	41.51600	42.82400					43.13000	4.852
20 PETN	+++++	+++++	+++++	+++++	+++++	+++++	+++++	+++++

TestAmerica West Sacramento

INITIAL CALIBRATION DATA

Start Cal Date : 10-JUN-2009 15:33  
 End Cal Date : 11-JUN-2009 13:47  
 Quant Method : ESTD  
 Origin : Disabled  
 Target Version : 4.14  
 Integrator : Falcon  
 Method file : \\Terastation\share\GCdata\LC10.I\06102009.B\8330AB.M  
 Last Edit : 11-Jun-2009 15:06 kenneyf  
 Curve Type : Average

Compound	5.000 Level 1	10.000 Level 2	20.000 Level 3	50.000 Level 4	100.000 Level 5	200.000 Level 6	RRF	% RSD
	500.000 Level 7	1000.000 Level 8						
\$ 1 3,4-Dinitrotoluene	+++++ 46.08000	50.10000 48.38600	46.40000	46.43000	45.87000	46.85500	47.16729	3.251

TestAmerica West Sacramento

INITIAL CALIBRATION DATA

Start Cal Date : 10-JUN-2009 16:23  
 End Cal Date : 11-JUN-2009 13:47  
 Quant Method : ESTD  
 Origin : Disabled  
 Target Version : 4.14  
 Integrator : Falcon  
 Method file : \\Terastation\share\GCdata\LC10.I\06102009.B\8330AB.M\83  
 Last Edit : 11-Jun-2009 15:14 kenneyf  
 Curve Type : Average

Calibration File Names:

Level 2: \\Terastation\share\GCdata\GCdata\LC10.I\06102009.B\A-000003.d\A-  
 Level 3: \\Terastation\share\GCdata\GCdata\LC10.I\06102009.B\A-000015.d\A-  
 Level 4: \\Terastation\share\GCdata\GCdata\LC10.I\06102009.B\A-000005.d\A-  
 Level 5: \\Terastation\share\GCdata\GCdata\LC10.I\06102009.B\A-000006.d\A-  
 Level 6: \\Terastation\share\GCdata\GCdata\LC10.I\06102009.B\A-000007.d\A-  
 Level 7: \\Terastation\share\GCdata\GCdata\LC10.I\06102009.B\A-000008.d\A-  
 Level 8: \\Terastation\share\GCdata\GCdata\LC10.I\06102009.B\A-000009.d\A-

Compound	10.000 Level 2	20.000 Level 3	50.000 Level 4	100.000 Level 5	200.000 Level 6	500.000 Level 7	1000.000 Level 8	RRF	% RSD
2 HMX	+++++	+++++	+++++	+++++	+++++	+++++		+++++	+++++
3 RDX	+++++	+++++	+++++	+++++	+++++	+++++		+++++	+++++
4 EGDN	+++++	+++++	+++++	+++++	+++++	+++++		+++++	+++++
5 Picric ACID	+++++	134	126	125	125	110	110	122	8.080
6 1,3,5-Trinitrobenzene	+++++	+++++	+++++	+++++	+++++	+++++		+++++	+++++
7 1,3-Dinitrobenzene	+++++	+++++	+++++	+++++	+++++	+++++		+++++	+++++
8 3,5-Dinitroaniline	+++++	+++++	+++++	+++++	+++++	+++++		+++++	+++++
9 TETRYL	+++++	+++++	+++++	+++++	+++++	+++++		+++++	+++++



TestAmerica West Sacramento

INITIAL CALIBRATION DATA

Start Cal Date : 10-JUN-2009 16:23  
 End Cal Date : 11-JUN-2009 13:47  
 Quant Method : ESTD  
 Origin : Disabled  
 Target Version : 4.14  
 Integrator : Falcon  
 Method file : \\Terastation\share\GCdata\LC10.I\06102009.B\8330AB.M\83  
 Last Edit : 11-Jun-2009 15:14 kenneyf  
 Curve Type : Average

Compound	10.000 Level 2	20.000 Level 3	50.000 Level 4	100.000 Level 5	200.000 Level 6	500.000 Level 7	RRF	% RSD
	1000.000 Level 8							
10 Nitrobenzene	++++	++++	++++	++++	++++	++++	++++	++++
11 Nitroglycerin	++++ 61.73900	61.25000	59.82000	60.58000	61.86500	59.75400	60.83467	1.527
12 2,4,6-Trinitrotoluene	++++	++++	++++	++++	++++	++++	++++	++++
13 4-AM-2,6-DNT	++++	++++	++++	++++	++++	++++	++++	++++
14 2-AM-4,6-DNT	++++	++++	++++	++++	++++	++++	++++	++++
15 2,6-Dinitrotoluene	++++	++++	++++	++++	++++	++++	++++	++++
16 2,4-Dinitrotoluene	++++	++++	++++	++++	++++	++++	++++	++++
17 2-Nitrotoluene	++++	++++	++++	++++	++++	++++	++++	++++
18 4-Nitrotoluene	++++	++++	++++	++++	++++	++++	++++	++++
19 3-Nitrotoluene	++++	++++	++++	++++	++++	++++	++++	++++
20 PETN	++++ 32.75200	32.65000	30.36000	31.91000	32.25000	32.38000	32.05033	2.748

TestAmerica West Sacramento

INITIAL CALIBRATION DATA

Start Cal Date : 10-JUN-2009 16:23  
 End Cal Date : 11-JUN-2009 13:47  
 Quant Method : ESTD  
 Origin : Disabled  
 Target Version : 4.14  
 Integrator : Falcon  
 Method file : \\Terastation\share\GCdata\LC10.I\06102009.B\8330AB.M\83  
 Last Edit : 11-Jun-2009 15:14 kenneyf  
 Curve Type : Average

Compound	10.000	20.000	50.000	100.000	200.000	500.000	RRF	% RSD
	Level 2	Level 3	Level 4	Level 5	Level 6	Level 7		
	1000.000							
	Level 8							
\$ 1 3,4-Dinitrotoluene	* +++++ 93.14200	90.40000	88.76000	88.28000	90.51000	88.03667	89.85478	2.142

\* Level 2 dropped due to poor integration  
 JLN  
 6/12/09

Chromatography Summary

Injection Date: 6/10/2009 23:09 Operator: fmk  
 DataFile: LC10.105102009.BA-000011.D Vial Num: 20  
 Instrument ID: LC10

Method 8330 Target Analyte Results

Method File: LC10.105102009.BA8330AB.M  
 Start Cal Date: 6/10/2009 15:33 End Cal Date: 6/11/2009 13:47

Sample: **ICV\_6 08GCSV0397**  
**100/200/100/100ng/mL**

Dilution Factor	Extract Volume	Sample Volume	Sample Weight
1X	0 mL	0 mL	0 g

Matrix: NONE SubList: CAL.sub SpikeList:  
 Samp. Info: ICV\_6 08GCSV0397 100/200/100/100ng/mL:2  
 Misc. Info: :6; ; ;3;CAL.sub; :0;1

Signal 1 UV 250-265								Signal 2 UV 358-205								
Compound Name	RT	Response	PPB	Spike Level	%D	Result	Flag	RT	Response	PPB	Spike Level	%D	Result	Flag	Limits(%)	Flag
<b>3,4-Dinitrotoluene</b>				<b>200</b>	<b>-100%</b>	<b>Fails</b>					<b>200</b>	<b>-100%</b>	<b>Fails</b>		<b>(±15)</b>	
HMX	5.29	26153	216.4000<	200	8%	Acceptable					200	-100%	Fails		(±15)	45
RDX	7.84	16084	206.5000<	200	3%	Acceptable					200	-100%	Fails		(±15)	45
Picric ACID	9.00	41754	504.4000 ✓	500	1%	Acceptable	9.00	61509	505.5000<	500	1%	Acceptable			(±15)	
1,3,5-Trinitrobenzene	10.20	29097	197.1000<	200	-1%	Acceptable					200	-100%	Fails		(±15)	45
1,3-Dinitrobenzene	13.25	27834	199.2000<	200	0%	Acceptable					200	-100%	Fails		(±15)	45
TETRYL	14.30	17776	191.6000<	200	-4%	Acceptable					200	-100%	Fails		(±15)	45
Nitrobenzene	15.16	12143	200.1000<	200	0%	Acceptable					200	-100%	Fails		(±15)	45
2,4,6-Trinitrotoluene	16.53	17352	197.8000<	200	-1%	Acceptable					200	-100%	Fails		(±15)	45
4-AM-2,6-DNT	17.51	11969	192.1000< ✓	200	-4%	Acceptable					200	-100%	Fails		(±15)	45
2-AM-4,6-DNT	18.57	13851	188.3000<	200	-6%	Acceptable					200	-100%	Fails		(±15)	45
2,6-Dinitrotoluene	20.28	10445	196.2000<	200	-2%	Acceptable					200	-100%	Fails		(±15)	45
2,4-Dinitrotoluene	21.00	16653	194.5000<	200	-3%	Acceptable					200	-100%	Fails		(±15)	45
2-Nitrotoluene	24.68	7069	196.7000<	200	-2%	Acceptable					200	-100%	Fails		(±15)	45
4-Nitrotoluene	26.50	8526	194.7000<	200	-3%	Acceptable					200	-100%	Fails		(±15)	45
3-Nitrotoluene	28.49	8451	195.9000< ✓	200	-2%	Acceptable					200	-100%	Fails		(±15)	45
<b>Nitroglycerin</b>				<b>200</b>	<b>-100%</b>	<b>Fails</b>		15.78	12548	206.3000< ✓	200	3%	Acceptable		(±15)	45
PE'IN				200	-100%	Fails		30.10	5630	181.9000< ✓	200	-9%	Acceptable		(±15)	45
3,5-Dinitroaniline	13.99	18460	205.2000<	200	3%	Acceptable					200	-100%	Fails		(±15)	45
<b>EGDN</b>				<b>200</b>	<b>-100%</b>	<b>Fails</b>					200	-100%	Fails		(±15)	

Notes: M = Manually Integrated Signals Differ by More Than 40%  
 D = Operator Disabled Result Signals Differ by More Than 50%  
 O = Over Calibration Range



Chromatography Summary

Injection Date: 7/9/2009 0:35 Operator: fhk  
 Data File: LC10 I07082009 BVA-000011.D Vial Num: 3  
 Instrument ID: LC10

Method 8330 Target Analyte Results

Sample: **STD\_5 09GCSV0053 .2K/.1/.1/.1**

Method File: LC10 I07082009 B8330AB.M  
 Start Cal Date: 6/10/2009 15:33 End Cal Date: 6/11/2009 13:47

Matrix: NONE SubList: CAL sub SpikeList:  
 Samp. Info: STD\_5 09GCSV0053 .2K/.1/.1/.1;  
 Misc. Info: .5; .; .3,CAL sub, .0,1

Dilution Factor	Extract Volume	Sample Volume	Sample Weight
1X	0 mL	0 mL	0 g

Signal 1 UV 250-265							Signal 2 UV 358-205									
Compound Name	RT	Response	PPB	Spike Level	%D	Result	Flag	RT	Response	PPB	Spike Level	%D	Result	Flag	Limits(%)	Flag
3,4-Dinitrotoluene	18.20	4739	100.5000<	100	1%	Acceptable		18.21	9053	104.8000	100	5%	Acceptable		(±15)	
HMX	5.29	12630	104.5000<	100	5%	Acceptable					100	-100%	Fails		(±15)	45
RDX	7.85	8029	103.1000<	100	3%	Acceptable					100	-100%	Fails		(±15)	45
Picric ACID	9.05	16777	202.7000	200	1%	Acceptable		9.06	24679	202.8000<	200	1%	Acceptable		(±15)	
1,3,5-Trinitrobenzene	10.20	15097	102.3000<	100	2%	Acceptable					100	-100%	Fails		(±15)	45
1,3-Dinitrobenzene	13.26	14436	103.3000<	100	3%	Acceptable					100	-100%	Fails		(±15)	45
TETRYL	14.34	9407	101.4000<	100	1%	Acceptable					100	-100%	Fails		(±15)	45
Nitrobenzene	15.18	6290	103.7000<	100	4%	Acceptable					100	-100%	Fails		(±15)	45
2,4,6-Trinitrotoluene	16.56	8976	102.3000<	100	2%	Acceptable					100	-100%	Fails		(±15)	45
4-AM-2,6-DNT	17.55	6365	102.1000<	100	2%	Acceptable					100	-100%	Fails		(±15)	45
2-AM-4,6-DNT	18.62	7472	101.6000<	100	2%	Acceptable					100	-100%	Fails		(±15)	45
2,6-Dinitrotoluene	20.35	5434	102.1000<	100	2%	Acceptable					100	-100%	Fails		(±15)	45
2,4-Dinitrotoluene	21.06	8697	101.6000<	100	2%	Acceptable					100	-100%	Fails		(±15)	45
2-Nitrotoluene	24.79	3631	101.0000<	100	1%	Acceptable					100	-100%	Fails		(±15)	45
4-Nitrotoluene	26.60	4405	100.6000<	100	1%	Acceptable					100	-100%	Fails		(±15)	45
3-Nitrotoluene	28.62	4379	101.5000<	100	2%	Acceptable					100	-100%	Fails		(±15)	45
Nitroglycerin				100	-100%	Fails		15.81	6320	103.9000<	100	4%	Acceptable		(±15)	45
PETN				100	-100%	Fails		30.24	3350	105.4000<	100	5%	Acceptable		(±15)	45
3,5-Dinitroaniline	14.01	9130	101.5000<	100	2%	Acceptable					100	-100%	Fails		(±15)	45
EGDN				100	-100%	Fails					100	-100%	Fails		(±15)	45

*NDF*

Notes: M = Manually Integrated Signals Differ by More Than 40%  
 D = Operator Disabled Result Signals Differ by More Than 50%  
 O = Over Calibration Range

Test America West Sacramento

Method 8330

Data file : \\Terastation\share\GCdata\LC10.I\07082009.B\A-000011.D  
 Lab Smp Id: STD\_5\_09GCSV0053\_2  
 Inj Date : 09-JUL-2009 00:35  
 Operator : fhk Inst ID: LC10.i  
 Smp Info : STD\_5\_09GCSV0053\_2K/.1/.1/.1/.1;2  
 Misc Info : ;5;-; ; ;3;CAL.sub; ;0;1  
 Comment : SOP SAC-LC-0009  
 Method : \\Terastation\share\GCdata\LC10.I\07082009.B\8330AB.m  
 Meth Date : 09-Jul-2009 01:21 tap Quant Type: AREA%  
 Cal Date : 11-JUN-2009 13:47 Cal File: A-000015.d  
 Als bottle: 3 Continuing Calibration Sample  
 Dil Factor: 1.00000  
 Integrator: Falcon+ Compound Sublist: CAL.sub  
 Target Version: 4.14  
 Processing Host: SACP307HPLC

RT	AREA	HEIGHT	HT/AREA	% HEIGHT	COMPOUNDS
5.289	81309	12630	0.155	8.64	2 HMX
7.849	92683	8029	0.087	5.49	3 RDX
8.599	1160	108	0.093	0.07	
9.052	241185	16777	0.070	11.58	5 Picric ACID
10.202	207248	15097	0.073	10.33	6 1,3,5-Trinitrobenze
12.785	119	30	0.252	0.02	
13.259	246389	14436	0.059	9.88	7 1,3-Dinitrobenzene
14.005	151716	9130	0.060	6.25	8 3,5-Dinitroaniline
14.335	168680	9407	0.056	6.43	9 TETRYL
15.179	121046	6290	0.052	4.30	10 Nitrobenzene
16.559	173239	8976	0.052	6.14	12 2,4,6-Trinitrotolue
17.552	134801	6365	0.047	4.35	13 4-AM-2,6-DNT
18.202	90656	4739	0.052	3.24	\$ 1 3,4-Dinitrotoluene
18.622	179304	7472	0.042	5.11	14 2-AM-4,6-DNT
20.349	124935	5434	0.043	3.71	15 2,6-Dinitrotoluene
21.062	214815	8697	0.040	5.95	16 2,4-Dinitrotoluene
22.369	505	45	0.089	0.03	
24.785	102226	3631	0.036	2.48	17 2-Nitrotoluene
26.599	132676	4405	0.033	3.01	18 4-Nitrotoluene
28.622	142181	4379	0.031	2.99	19 3-Nitrotoluene
	2606872	146077		100.000	

Total unknown % height = 0.1200

Chromatography Summary

Injection Date: 7/9/2009 5:38 Operator: fhk  
 Data File: LC10 I07082009 BVA-000017.D Vial Num: 4  
 Instrument ID: LC10

Method 8330 Target Analyte Results

Sample: **STD\_6 09GCSV0054 .4K/././././.**

Method File: LC10 I07082009 B8330AB.M  
 Start Cal Date: 6/10/2009 15:33 End Cal Date: 6/11/2009 13:47

Dilution Factor	Extract Volume	Sample Volume	Sample Weight
1X	0 mL	0 mL	0 g

Matrix: NONE SubList: CAL sub SpikeList:  
 Samp. Info: STD\_6 09GCSV0054 4K/././././.  
 Misc. Info: .6., .3;CAL sub; ;0;1

Compound Name	RT	Signal 1 UV 250-265					Signal 2 UV 358-205					Limits(%)	Flag	
		Response	PPB	Spike Level	%D	Result	RT	Response	PPB	Spike Level	%D			Result
3,4-Dinitrotoluene	18.21	9682	205.3000<	200	3%	Acceptable	18.21	18470	213.6000	200	7%	Acceptable	(±15)	
HMX	5.29	25614	212.0000<	200	6%	Acceptable				200	-100%	Fails	(±15)	45
RDX	7.85	16441	211.1000<	200	6%	Acceptable				200	-100%	Fails	(±15)	45
Picric ACID	9.03	42453	512.9000	500	3%	Acceptable	9.03	62382	512.7000<	500	3%	Acceptable	(±15)	
1,3,5-Trinitrobenzene	10.20	30599	207.3000<	200	4%	Acceptable				200	-100%	Fails	(±15)	45
1,3-Dinitrobenzene	13.26	29254	209.4000<	200	5%	Acceptable				200	-100%	Fails	(±15)	45
TETRYL	14.34	19071	205.5000<	200	3%	Acceptable				200	-100%	Fails	(±15)	45
Nitrobenzene	15.18	12804	211.0000<	200	6%	Acceptable				200	-100%	Fails	(±15)	45
2,4,6-Trinitrotoluene	16.56	18126	206.6000<	200	3%	Acceptable				200	-100%	Fails	(±15)	45
4-AM-2,6-DNT	17.56	12935	207.6000<	200	4%	Acceptable				200	-100%	Fails	(±15)	45
2-AM-4,6-DNT	18.62	15164	206.1000<	200	3%	Acceptable				200	-100%	Fails	(±15)	45
2,6-Dinitrotoluene	20.35	10941	205.5000<	200	3%	Acceptable				200	-100%	Fails	(±15)	45
2,4-Dinitrotoluene	21.07	17582	205.3000<	200	3%	Acceptable				200	-100%	Fails	(±15)	45
2-Nitrotoluene	24.78	7317	203.6000<	200	2%	Acceptable				200	-100%	Fails	(±15)	45
4-Nitrotoluene	26.60	8872	202.6000<	200	1%	Acceptable				200	-100%	Fails	(±15)	45
3-Nitrotoluene	28.62	8820	204.5000<	200	2%	Acceptable				200	-100%	Fails	(±15)	45
Nitroglycerin				200	-100%	Fails	15.81	12540	206.1000<	200	3%	Acceptable	(±15)	45
PETN				200	-100%	Fails	30.24	6614	207.0000<	200	4%	Acceptable	(±15)	45
3,5-Dinitroaniline	14.01	18615	206.9000<	200	3%	Acceptable				200	-100%	Fails	(±15)	45
EGDN				200	-100%	Fails				200	-100%	Fails	(±15)	

Notes M = Manually Integrated Signals Differ by More Than 40%  
 D = Operator Disabled Result Signals Differ by More Than 50%  
 O = Over Calibration Range





Chromatography Summary

Injection Date: 7/11/2009 4:29 Operator: fhk  
 DataFile: LC10.N07102009.BVA-000021.D Vial Num: 4  
 Instrument ID: LC10

Method 8330 Target Analyte Results

Sample: STD\_5 09GCSV0238 2K/1/1/1/1

Method File: LC10.N07102009.BVA8330AB.M  
 Start Cal Date: 5/10/2009 15:33 End Cal Date: 6/11/2009 13:47

Matrix: NONE SubList: CAL.sub SpikeList:  
 Samp. Info: STD\_5 09GCSV0238 2K/1/1/1/1,2  
 Misc. Info: ;5; ; ;3;CAL sub; 0;1

Dilution Factor	Extract Volume	Sample Volume	Sample Weight
1X	0 mL	0 mL	0 g

Signal 1 UV 250-265								Signal 2 UV 358-205								
Compound Name	RT	Response	PPB	Spike Level	%D	Result	Flag	RT	Response	PPB	Spike Level	%D	Result	Flag	Limits(%)	Flag
3,4-Dinitrotoluene	18.23	4761	100.9000<	100	1%	Acceptable		18.22	9142	105.8000	100	6%	Acceptable		(±15)	
HMX	5.29	12731	105.4000<	100	5%	Acceptable					100	-100%	Fails		(±15)	45
RDX	7.85	8006	102.8000<	100	3%	Acceptable					100	-100%	Fails		(±15)	45
Picric ACID	9.07	17082	206.4000	200	3%	Acceptable		9.07	25079	206.1000<	200	3%	Acceptable		(±15)	
1,3,5-Trinitrobenzene	10.21	14836	100.5000<	100	1%	Acceptable					100	-100%	Fails		(±15)	45
1,3-Dinitrobenzene	13.27	14248	102.0000<	100	2%	Acceptable					100	-100%	Fails		(±15)	45
TETRYL	14.35	9417	101.5000<	100	2%	Acceptable					100	-100%	Fails		(±15)	45
Nitrobenzene	15.19	6384	105.2000<	100	5%	Acceptable					100	-100%	Fails		(±15)	45
2,4,6-Trinitrotoluene	16.57	9006	102.7000<	100	3%	Acceptable					100	-100%	Fails		(±15)	45
4-AM-2,6-DNT	17.57	6469	103.8000<	100	4%	Acceptable					100	-100%	Fails		(±15)	45
2-AM-4,6-DNT	18.63	7621	103.6000<	100	4%	Acceptable					100	-100%	Fails		(±15)	45
2,6-Dinitrotoluene	20.36	5436	102.1000<	100	2%	Acceptable					100	-100%	Fails		(±15)	45
2,4-Dinitrotoluene	21.08	8612	100.6000<	100	1%	Acceptable					100	-100%	Fails		(±15)	45
2-Nitrotoluene	24.80	3719	103.5000<	100	4%	Acceptable					100	-100%	Fails		(±15)	45
4-Nitrotoluene	26.63	4483	102.4000<	100	2%	Acceptable					100	-100%	Fails		(±15)	45
3-Nitrotoluene	28.64	4429	102.7000<	100	3%	Acceptable					100	-100%	Fails		(±15)	45
Nitroglycerin				100	-100%	Fails		15.83	6116	100.5000<	100	1%	Acceptable		(±15)	45
PETN				100	-100%	Fails		30.27	3354	104.6000<	100	5%	Acceptable		(±15)	45
3,5-Dinitroaniline	14.01	9259	102.9000<	100	3%	Acceptable					100	-100%	Fails		(±15)	45
EGDN				100	-100%	Fails					100	-100%	Fails		(±15)	

Notes M = Manually Integrated Signals Differ by More Than 40%  
 D = Operator Disabled Result Signals Differ by More Than 50%  
 O = Over Calibration Range

Chromatography Summary

Injection Date: 7/11/2009 12:55 Operator: fhk  
 DataFile: LC10 107102009 BVA-000031.D Vial Num: 5  
 Instrument ID: .C10

Method 8330 Target Analyte Results

Sample: STD\_5 09GCSV0238 .2K/1/1/1/1

Method File: LC10 107102009 BVA8330AB.M  
 Start Cal Date: 6/10/2009 15:33 End Cal Date: 6/11/2009 13:47

Matrix: NONE SubList: CAL.sub SpikeList:  
 Samp. Info: STD\_5 09GCSV0238 .2K/1/1/1/1.2  
 Misc. Info: .5, ., .3, CAL.sub, .0, 1

Dilution Factor	Extract Volume	Sample Volume	Sample Weight
1X	0 mL	0 mL	0 g

Compound Name	RT	Signal 1 UV 250-265					Signal 2 UV 358-205					Limits(%)	Flag	
		Response	PPB	Spike Level	%D	Result	RT	Response	PPB	Spike Level	%D			Result
3,4-Dinitrotoluene	18.24	4768	101.1000<	100	1%	Acceptable	18.25	9178	106.2000	100	6%	Acceptable	(±15)	
HMX	5.29	12737	105.4000<	100	5%	Acceptable				100	-100%	Fails	(±15)	45
RDX	7.86	8013	102.9000<	100	3%	Acceptable				100	-100%	Fails	(±15)	45
Picric ACID	9.07	17075	206.3000	200	3%	Acceptable	9.08	25051	205.9000<	200	3%	Acceptable	(±15)	
1,3,5-Trinitrobenzene	10.21	14849	100.6000<	100	1%	Acceptable				100	-100%	Fails	(±15)	45
1,3-Dinitrobenzene	13.28	14249	102.0000<	100	2%	Acceptable				100	-100%	Fails	(±15)	45
TETRYL	14.37	9416	101.5000<	100	2%	Acceptable				100	-100%	Fails	(±15)	45
Nitrobenzene	15.21	6360	104.8000<	100	5%	Acceptable				100	-100%	Fails	(±15)	45
2,4,6-Trinitrotoluene	16.59	8994	102.5000<	100	3%	Acceptable				100	-100%	Fails	(±15)	45
4-AM-2,6-DNT	17.60	6456	103.6000<	100	4%	Acceptable				100	-100%	Fails	(±15)	45
2-AM-4,6-DNT	18.66	7636	103.8000<	100	4%	Acceptable				100	-100%	Fails	(±15)	45
2,6-Dinitrotoluene	20.39	5431	102.0000<	100	2%	Acceptable				100	-100%	Fails	(±15)	45
2,4-Dinitrotoluene	21.10	8609	100.5000<	100	1%	Acceptable				100	-100%	Fails	(±15)	45
2-Nitrotoluene	24.83	3699	102.9000<	100	3%	Acceptable				100	-100%	Fails	(±15)	45
4-Nitrotoluene	26.66	4472	102.1000<	100	2%	Acceptable				100	-100%	Fails	(±15)	45
3-Nitrotoluene	28.68	4415	102.4000<	100	2%	Acceptable				100	-100%	Fails	(±15)	45
Nitroglycerin				100	-100%	Fails	15.85	6140	100.9000<	100	1%	Acceptable	(±15)	45
PEFN				100	-100%	Fails	30.33	3344	104.3000<	100	4%	Acceptable	(±15)	45
3,5-Dinitroaniline	14.03	9257	102.9000<	100	3%	Acceptable				100	-100%	Fails	(±15)	45
EGDN				100	-100%	Fails				100	-100%	Fails	(±15)	

Notes: M = Manually Integrated Signals Differ by More Than 40%  
 D = Operator Disabled Result Signals Differ by More Than 50%  
 O = Over Calibration Range



TestAmerica West Sacramento

INITIAL CALIBRATION DATA

Start Cal Date : 09-JUN-2009 20:53  
 End Cal Date : 10-JUN-2009 04:20  
 Quant Method : ESTD  
 Origin : Disabled  
 Target Version : 4.14  
 Integrator : HP Genie  
 Method file : \\Terastation\share\GCdata\LC9.I\06092009.B\8330CNAB.M  
 Last Edit : 10-Jun-2009 10:24 kenneyf  
 Curve Type : Average

Calibration File Names:

- Level 1: \\Terastation\share\GCdata\GCdata\LC9.I\06092009.B\C-000004.d
- Level 2: \\Terastation\share\GCdata\GCdata\LC9.I\06092009.B\C-000005.d
- Level 3: \\Terastation\share\GCdata\GCdata\LC9.I\06092009.B\C-000006.d
- Level 4: \\Terastation\share\GCdata\GCdata\LC9.I\06092009.B\C-000007.d
- Level 5: \\Terastation\share\GCdata\GCdata\LC9.I\06092009.B\C-000008.d
- Level 6: \\Terastation\share\GCdata\GCdata\LC9.I\06092009.B\C-000009.d
- Level 7: \\Terastation\share\GCdata\GCdata\LC9.I\06092009.B\C-000010.d
- Level 8: \\Terastation\share\GCdata\GCdata\LC9.I\06092009.B\C-000011.d

Compound	5.000	10.000	20.000	50.000	100.000	200.000	RRF	% RSD
	Level 1	Level 2	Level 3	Level 4	Level 5	Level 6		
2 HMX	49.60000 37.41800	43.70000 39.22000	40.45000	40.46000	39.51000	39.31500	41.20913	9.288
3 RDX	45.40000 38.21800	43.70000 42.45200	42.75000	43.30000	42.16000	42.58500	42.55812	4.852
4 EGDN	++++ ++++	++++ ++++	++++	++++	++++	++++	++++	++++ <-
5 Picric ACID	++++ ++++	++++ ++++	++++	++++	++++	++++	++++	++++ <-
6 1,3,5-Trinitrobenzene	66.80000 63.46600	67.20000 67.27500	67.60000	68.82000	67.02000	67.61500	66.97450	2.306
7 1,3-Dinitrobenzene	95.20000 79.37000	88.70000 88.38800	87.25000	89.40000	87.58000	88.56500	88.05662	4.893
8 3,5-Dinitroaniline	71.20000 60.97200	68.50000 65.89900	68.65000	68.54000	66.12000	66.97500	67.10700	4.478
9 TETRYL	93.20000 85.04400	86.50000 86.33600	85.70000	86.92000	86.05000	86.26500	87.00187	2.950

TestAmerica West Sacramento

INITIAL CALIBRATION DATA

Start Cal Date : 09-JUN-2009 20:53  
 End Cal Date : 10-JUN-2009 04:20  
 Quant Method : ESTD  
 Origin : Disabled  
 Target Version : 4.14  
 Integrator : HP Genie  
 Method file : \\Terastation\share\GCdata\LC9.I\06092009.B\8330CNAB.M  
 Last Edit : 10-Jun-2009 10:24 kenneyf  
 Curve Type : Average

Compound	5.000	10.000	20.000	50.000	100.000	200.000	RRF	% RSD
	Level 1	Level 2	Level 3	Level 4	Level 5	Level 6		
	500.000	1000.000						
	Level 7	Level 8						
10 Nitrobenzene	39.40000 35.35200	42.00000 40.30500	38.65000	40.64000	39.79000	40.66500	39.60025	5.005
11 Nitroglycerin	+++++	+++++	+++++	+++++	+++++	+++++	+++++	+++++
12 2,4,6-Trinitrotoluene	104 71.97000	91.00000 73.60800	81.00000	77.16000	74.32000	73.90500	80.89538	13.892
13 4-AM-2,6-DNT	68.80000 64.80800	69.80000 65.14700	68.05000	67.74000	65.37000	65.85500	66.94625	2.818
14 2-AM-4,6-DNT	80.20000 69.79400	76.80000 72.19800	75.45000	74.50000	72.25000	73.07500	74.28838	4.343
15 2,6-Dinitrotoluene	54.80000 45.69200	52.40000 49.41100	49.70000	49.88000	49.11000	49.44000	50.05412	5.280
16 2,4-Dinitrotoluene	83.00000 73.10400	82.40000 78.73900	79.60000	80.22000	78.86000	79.32000	79.40538	3.779
17 2-Nitrotoluene	24.50000 22.18600	25.30000 24.28200	23.52500	24.39000	24.32500	24.51250	24.12756	3.813
18 4-Nitrotoluene	+++++	+++++	+++++	+++++	+++++	+++++	+++++	+++++
19 3-Nitrotoluene	35.00000 28.82000	32.50000 30.96400	30.55000	31.42000	30.76000	31.17500	31.39862	5.672
20 PETN	+++++	+++++	+++++	+++++	+++++	+++++	+++++	+++++

TestAmerica West Sacramento

INITIAL CALIBRATION DATA

Start Cal Date : 09-JUN-2009 20:53  
 End Cal Date : 10-JUN-2009 04:20  
 Quant Method : ESTD  
 Origin : Disabled  
 Target Version : 4.14  
 Integrator : HP Genie  
 Method file : \\Terastation\share\GCdata\LC9.I\06092009.B\8330CNAB.M  
 Last Edit : 10-Jun-2009 10:24 kenneyf  
 Curve Type : Average

Compound	5.000 Level 1	10.000 Level 2	20.000 Level 3	50.000 Level 4	100.000 Level 5	200.000 Level 6	RRF	% RSD
	500.000 Level 7	1000.000 Level 8						
\$ 1 3,4-Dinitrotoluene	++++	46.40000	40.85000	39.44000	38.04000	38.34500		
	35.87333	38.62800					39.65376	8.405

TestAmerica West Sacramento

INITIAL CALIBRATION DATA

Start Cal Date : 09-JUN-2009 21:56  
 End Cal Date : 10-JUN-2009 04:20  
 Quant Method : ESTD  
 Origin : Disabled  
 Target Version : 4.14  
 Integrator : HP Genie  
 Method file : \\Terastation\share\GCdata\LC9.I\06092009.B\8330CNAB.M\8  
 Last Edit : 10-Jun-2009 10:25 kenneyf  
 Curve Type : Average

Calibration File Names:

Level 2: \\Terastation\share\GCdata\GCdata\LC9.I\06092009.B\C-000005.d\C-0  
 Level 3: \\Terastation\share\GCdata\GCdata\LC9.I\06092009.B\C-000006.d\C-0  
 Level 4: \\Terastation\share\GCdata\GCdata\LC9.I\06092009.B\C-000007.d\C-0  
 Level 5: \\Terastation\share\GCdata\GCdata\LC9.I\06092009.B\C-000008.d\C-0  
 Level 6: \\Terastation\share\GCdata\GCdata\LC9.I\06092009.B\C-000009.d\C-0  
 Level 7: \\Terastation\share\GCdata\GCdata\LC9.I\06092009.B\C-000010.d\C-0  
 Level 8: \\Terastation\share\GCdata\GCdata\LC9.I\06092009.B\C-000011.d\C-0

Compound	10.000 Level 2	20.000 Level 3	50.000 Level 4	100.000 Level 5	200.000 Level 6	500.000 Level 7	RRF	% RSD
2 HMX	++++	++++	++++	++++	++++	++++	++++	++++
3 RDX	++++	++++	++++	++++	++++	++++	++++	++++
4 EGDN	++++ 42.86600	39.00000	42.36000	44.25000	44.94000	40.36600	42.29700	5.363
5 Picric ACID	++++	++++	++++	++++	++++	++++	++++	++++ <-
6 1,3,5-Trinitrobenzene	++++	++++	++++	++++	++++	++++	++++	++++
7 1,3-Dinitrobenzene	++++	++++	++++	++++	++++	++++	++++	++++
8 3,5-Dinitroaniline	++++	++++	++++	++++	++++	++++	++++	++++
9 TETRYL	++++	++++	++++	++++	++++	++++	++++	++++

TestAmerica West Sacramento

INITIAL CALIBRATION DATA

Start Cal Date : 09-JUN-2009 21:56  
 End Cal Date : 10-JUN-2009 04:20  
 Quant Method : ESTD  
 Origin : Disabled  
 Target Version : 4.14  
 Integrator : HP Genie  
 Method file : \\Terastation\share\GCdata\LC9.I\06092009.B\8330CNAB.M\8  
 Last Edit : 10-Jun-2009 10:25 kenneyf  
 Curve Type : Average

Compound	10.000 Level 2	20.000 Level 3	50.000 Level 4	100.000 Level 5	200.000 Level 6	500.000 Level 7	RRP	% RSD
	1000.000 Level 8							
10 Nitrobenzene	+++++	+++++	+++++	+++++	+++++	+++++	+++++	+++++
11 Nitroglycerin	+++++ 57.30100	55.25000	55.16000	56.31000	57.02000	56.37800	56.23650	1.571
12 2,4,6-Trinitrotoluene	+++++	+++++	+++++	+++++	+++++	+++++	+++++	+++++
13 4-AM-2,6-DNT	+++++	+++++	+++++	+++++	+++++	+++++	+++++	+++++
14 2-AM-4,6-DNT	+++++	+++++	+++++	+++++	+++++	+++++	+++++	+++++
15 2,6-Dinitrotoluene	+++++	+++++	+++++	+++++	+++++	+++++	+++++	+++++
16 2,4-Dinitrotoluene	+++++	+++++	+++++	+++++	+++++	+++++	+++++	+++++
17 2-Nitrotoluene	+++++	+++++	+++++	+++++	+++++	+++++	+++++	+++++
18 4-Nitrotoluene	+++++	+++++	+++++	+++++	+++++	+++++	+++++	+++++
19 3-Nitrotoluene	+++++	+++++	+++++	+++++	+++++	+++++	+++++	+++++
20 PETN	+++++ 105	104	101	102	104	104	103	1.208



TestAmerica West Sacramento

INITIAL CALIBRATION DATA

Start Cal Date : 09-JUN-2009 21:56  
 End Cal Date : 10-JUN-2009 04:20  
 Quant Method : ESTD  
 Origin : Disabled  
 Target Version : 4.14  
 Integrator : HP Genie  
 Method file : \\Terastation\share\GCdata\LC9.I\06092009.B\8330CNAB.M\8  
 Last Edit : 10-Jun-2009 10:25 kenneyf  
 Curve Type : Average

Compound	10.000 Level 2	20.000 Level 3	50.000 Level 4	100.000 Level 5	200.000 Level 6	500.000 Level 7	RRP	% RSD
	1000.000 Level 8							
\$ 1,3,4-Dinitrotoluene	+++++	+++++	+++++	+++++	+++++	+++++	+++++	+++++

Chromatography Summary

Injection Date: 6/10/2009 6:29 Operator: fhk  
 DataFile: LC9 I06092009.BIC-000013.D Vial Num: 20  
 Instrument ID: LC9

Method 8330 Target Analyte Results

Sample: **ICV\_6 08GCSV0397**  
**200/500/200/200/200**

Method File: LC9 I06092009.BI8330CNA8.M  
 Start Cal Date: 6/9/2009 20:53 End Cal Date: 6/10/2009 4:20

Matrix: NONE SubList: CAL.sub SpikeList:  
 Samp. Info: ICV\_6 08GCSV0397 200/500/200/200/200;2  
 Misc. Info: , 6, , , ; 3; CAL.sub, . 0, 1

Dilution Factor	Extract Volume	Sample Volume	Sample Weight
1X	0 mL	0 mL	0 g

Signal 1 UV 250-265								Signal 2 UV 358-205								
Compound Name	RT	Response	PPB	Spike Level	%D	Result	Flag	RT	Response	PPB	Spike Level	%D	Result	Flag	Limits(%)	Flag
3,4-Dinitrotoluene	31.18	259	6.5320<	200	97%	Fails	Not in list				200	-100%	Fails		(±15)	45
HMX	35.11	8323	202.0000<	200	1%	Acceptable					200	-100%	Fails		(±15)	45
RDX	25.60	8556	201.0000<	200	1%	Acceptable					200	-100%	Fails		(±15)	45
Picric ACID				500	-100%	Fails					500	-100%	Fails		(±15)	
1,3,5-Trinitrobenzene	21.78	13318	198.8000<	200	-1%	Acceptable					200	-100%	Fails		(±15)	45
1,3-Dinitrobenzene	19.50	17403	197.6000<	200	-1%	Acceptable					200	-100%	Fails		(±15)	45
TETRYL	40.80	16134	185.4000<	200	-7%	Acceptable					200	-100%	Fails		(±15)	45
Nitrobenzene	16.81	7967	201.2000<	200	1%	Acceptable					200	-100%	Fails		(±15)	45
2,4,6-Trinitrotoluene	32.92	14562	180.0000<	200	-10%	Acceptable					200	-100%	Fails		(±15)	45
4-AM-2,6-DNT	29.55	12516	187.0000<	200	-7%	Acceptable					200	-100%	Fails		(±15)	45
2-AM-4,6-DNT	29.21	14216	191.4000<	200	-4%	Acceptable					200	-100%	Fails		(±15)	45
2,6-Dinitrotoluene	27.26	9755	194.9000<	200	-3%	Acceptable					200	-100%	Fails		(±15)	45
2,4-Dinitrotoluene	26.42	15486	195.0000<	200	-3%	Acceptable					200	-100%	Fails		(±15)	45
2-Nitrotoluene	22.83	9684	401.4000<	400	0%	Acceptable					400	-100%	Fails		(±15)	45
4-Nitrotoluene				400	-100%	Fails					400	-100%	Fails		(±15)	
3-Nitrotoluene	23.35	6108	194.5000<	200	-3%	Acceptable					200	-100%	Fails		(±15)	45
Nitroglycerin				200	-100%	Fails		38.67	11785	209.6000<	200	5%	Acceptable		(±15)	45
PETN				200	-100%	Fails		49.13	19084	184.7000<	200	-8%	Acceptable		(±15)	45
3,5-Dinitroaniline	25.06	13705	204.2000<	200	2%	Acceptable					200	-100%	Fails		(±15)	45
EGDN				200	-100%	Fails		18.01	9041	213.8000<	200	7%	Acceptable		(±15)	45

Notes: M = Manually Integrated Signals Differ by More Than 40%  
 D = Operator Disabled Result Signals Differ by More Than 50%  
 O = Over Calibration Range

TestAmerica West Sacramento  
GC/LC INSTRUMENT LOG

Inst ID: LC9 Batch ID: 07102009A  
Method : Method 8330 Test : SOP WS-LC-0009  
ICAL Date: See Calibration Report

Date	Time	Operator	Sample ID	File ID	Vol or Wt	Extract Vol	Diln	Comments
10-JUL-2009	17:28	fhk	Primer	C-000001	0 g	0 mL	1	
10-JUL-2009	18:30	fhk	Primer	C-000002	0 g	0 mL	1	
10-JUL-2009	19:32	fhk	STD_5 09GCSV0238 .2K/.1/ 1/ 1/	C-000003	0 g	0 mL	1	
10-JUL-2009	20:34	fhk	LF1EN1AA 9183251 G9G020000-251	C-000004	1000 mL	20 mL	1	
10-JUL-2009	21:36	fhk	LFXXW1AC 9183251 G9G010225-1 1	C-000005	1021.57 mL	20 mL	1	
10-JUL-2009	22:39	fhk	LPX111AC 9183251 G9G010225-3 1	C-000006	1024.15 mL	20 mL	1	
10-JUL-2009	23:41	fhk	LPX121AC 9183251 G9G010225-4 1	C-000007	1019.2 mL	20 mL	1	
11-JUL-2009	00:43	fhk	LPX151AC 9183251 G9G010225-5 1	C-000008	1021.92 mL	20 mL	1	
11-JUL-2009	01:45	fhk	LF7A31AA 9189394 G9G080000-394	C-000009	10 g	80 mL	1	
11-JUL-2009	02:48	fhk	LPX4Q1AA 9189394 G9G010246-1 1	C-000010	10.01 g	80 mL	1	
11-JUL-2009	03:50	fhk	LPX4R1AA 9189394 G9G010246-2 1	C-000011	9.06 g	80 mL	1	
11-JUL-2009	04:52	fhk	LPX4T1AA 9189394 G9G010246-3 1	C-000012	10.01 g	80 mL	1	
11-JUL-2009	05:54	fhk	STD_6 09GCSV0054 .5K/.2/.2/ 2/	C-000013	0 g	0 mL	1	
11-JUL-2009	06:57	fhk	LF3161AA 9187206 G9G060000-206	C-000014	1000 mL	20 mL	1	
11-JUL-2009	07:59	fhk	LF3H41AC 9187206 G9G030156-1 1	C-000015	1000 mL	20 mL	1	1007.3 mL
11-JUL-2009	09:01	fhk	LF3H51AC 9187206 G9G030156-2 1	C-000016	1000 mL	20 mL	1	996.15
11-JUL-2009	10:03	fhk	LF3H61AC 9187206 G9G030156-3 5	C-000017	1000 mL	20 mL	5	1005.3 ↓
11-JUL-2009	11:06	fhk	LF5WA1AA 9188431 G9G070000-431	C-000018	2 g	40 mL	1	
11-JUL-2009	12:08	fhk	LPXP31AA 9188431 A9G010193-1 1	C-000019	2 g	40 mL	1	
11-JUL-2009	13:10	fhk	LPXQQ1AA 9188431 A9G010193-2 1	C-000020	2.02 g	40 mL	1	
11-JUL-2009	14:12	fhk	LPW791AAB 9182192 MB 1X	C-000021	1000 mL	20 mL	1	
11-JUL-2009	15:14	fhk	LPV9M1AC 9182192 G9F300242-5 1	C-000022	1016.72 mL	20 mL	1	
11-JUL-2009	16:16	fhk	STD_5 09GCSV0238 .2K/.1/.1/	C-000023	0 g	0 mL	1	

Chromatography Summary

Injection Date: 7/11/2009 5:54 Operator: fhk  
 DataFile: LC9 I07102009A BVC-000013.D Vial Num: 3  
 Instrument ID: LC9

Method 8330 Target Analyte Results

Sample: STD\_6 09GCSV0054 .5K/.2/.2/.2

Method File: LC9 I07102009A BV8330CNAB.M  
 Start Cal Date: 6/9/2009 20:53 End Cal Date: 6/10/2009 4:20

Matrix: NONE SubList: CAL sub SpikeList:  
 Samp. Info: STD\_6 09GCSV0054 .5K/.2/.2/.2  
 Misc. Info: , 6, . . . 3, CAL sub.; 0; 1; 0

Dilution Factor	Extract Volume	Sample Volume	Sample Weight
1X	0 mL	0 mL	0 g

Signal 1 UV 250-265								Signal 2 UV 358-205								
Compound Name	RT	Response	PPB	Spike Level	%D	Result	Flag	RT	Response	PPB	Spike Level	%D	Result	Flag	Limits(%)	Flag
3,4-Dinitrotoluene	30.69	8487	214.0000<	200	7%	Acceptable					200	-100%	Fails		(±15)	45
HMX	34.67	7594	184.3000<	200	-8%	Acceptable					200	-100%	Fails		(±15)	45
RDX	35.16	8673	203.8000<	200	2%	Acceptable					200	-100%	Fails		(±15)	45
<b>Picric ACID</b>				<b>500</b>	<b>-100%</b>	<b>Fails</b>					<b>500</b>	<b>-100%</b>	<b>Fails</b>		<b>(±15)</b>	
1,3,5-Trinitrobenzene	21.54	13992	208.9000<	200	4%	Acceptable					200	-100%	Fails		(±15)	45
1,4-Dinitrobenzene	19.18	18240	207.1000<	200	4%	Acceptable					200	-100%	Fails		(±15)	45
TETRYL	40.38	18246	209.7000<	200	5%	Acceptable					200	-100%	Fails		(±15)	45
Nitrobenzene	16.47	8291	209.4000<	200	5%	Acceptable					200	-100%	Fails		(±15)	45
2,4,6-Trinitrotoluene	32.59	16243	200.8000<	200	0%	Acceptable					200	-100%	Fails		(±15)	45
4-AM-2,6-DNT	29.08	13363	199.6000<	200	0%	Acceptable					200	-100%	Fails		(±15)	45
2-AM-4,6-DNT	28.74	14987	201.7000<	200	1%	Acceptable					200	-100%	Fails		(±15)	45
2,6-Dinitrotoluene	26.81	10167	203.1000<	200	2%	Acceptable					200	-100%	Fails		(±15)	45
2,4-Dinitrotoluene	25.99	16528	208.1000<	200	4%	Acceptable					200	-100%	Fails		(±15)	45
2-Nitrotoluene	22.38	10164	421.3000<	400	5%	Acceptable					400	-100%	Fails		(±15)	45
<b>4-Nitrotoluene</b>				<b>400</b>	<b>-100%</b>	<b>Fails</b>					<b>400</b>	<b>-100%</b>	<b>Fails</b>		<b>(±15)</b>	
3-Nitrotoluene	22.90	6323	201.4000<	200	1%	Acceptable					200	-100%	Fails		(±15)	45
<b>Nitroglycerin</b>				<b>200</b>	<b>-100%</b>	<b>Fails</b>		38.25	11803	209.9000<	200	5%	Acceptable		(±15)	45
<b>PETN</b>				<b>200</b>	<b>-100%</b>	<b>Fails</b>		48.78	21503	208.1000<	200	4%	Acceptable		(±15)	45
3,5-Dinitroaniline	24.65	13757	205.0000<	200	3%	Acceptable					200	-100%	Fails		(±15)	45
<b>EGDN</b>				<b>200</b>	<b>-100%</b>	<b>Fails</b>		17.80	9381	221.8000	200	11%	Acceptable		(±15)	45

Notes: M = Manually Integrated Signals Differ by More Than 40%  
 D = Operator Disabled Result Signals Differ by More Than 50%  
 O = Over Calibration Range

Chromatography Summary

Injection Date: 7/11/2009 16:16 Operator: fhk  
 DataFile: LC9\107102009A B\C-000023.D Vial Num: 4  
 Instrument ID: LC9

Method 8330 Target Analyte Results

Method File: LC9\107102009A B\8330CNAB.M  
 Start Cal Date: 6/9/2009 20:53 End Cal Date: 6/10/2009 4:20

Sample : **STD\_5 09GCSV0238 .2K/1/1/1/1**

Dilution Factor	Extract Volume	Sample Volume	Sample Weight
1X	0 mL	0 mL	0 g

Matrix: NONE SubList: CAL.sub SpikeList:  
 Samp. Info: STD\_5 09GCSV0238 2K/1/1/1/1,2  
 Misc. Info: .5. . ; 3, CAL.sub; . 0, 1, 0

Signal 1 UV 250-265								Signal 2 UV 358-205								
Compound Name	RT	Response	PPB	Spike Level	%D	Result	Flag	RT	Response	PPB	Spike Level	%D	Result	Flag	Limits(%)	Flag
3,4-Dinitrotoluene	30.65	4159	101.9000<	100	5%	Acceptable					100	-100%	Fails		(±15)	45
HMX	34.66	3754	91.1000<	100	-9%	Acceptable					100	-100%	Fails		(±15)	45
RDX	25.13	4283	100.6000<	100	1%	Acceptable					100	-100%	Fails		(±15)	45
<b>Picric ACID</b>				<b>200</b>	<b>-100%</b>	<b>Fails</b>					<b>200</b>	<b>-100%</b>	<b>Fails</b>		(±15)	
1,3,5-Trinitrobenzene	21.50	6847	102.2000<	100	2%	Acceptable					100	-100%	Fails		(±15)	45
1,3-Dinitrobenzene	19.16	8905	101.1000<	100	1%	Acceptable					100	-100%	Fails		(±15)	45
TETRYL	40.39	8996	103.4000<	100	3%	Acceptable					100	-100%	Fails		(±15)	45
Nitrobenzene	16.46	4129	104.3000<	100	4%	Acceptable					100	-100%	Fails		(±15)	45
2,4,6-Trinitrotoluene	32.52	7906	97.7300<	100	-2%	Acceptable					100	-100%	Fails		(±15)	45
4-AM-2,6-DNT	29.04	6874	102.7000<	100	3%	Acceptable					100	-100%	Fails		(±15)	45
2-AM-4,6-DNT	28.71	7546	101.6000<	100	2%	Acceptable					100	-100%	Fails		(±15)	45
2,6-Dinitrotoluene	26.77	5040	100.7000<	100	1%	Acceptable					100	-100%	Fails		(±15)	45
2,4-Dinitrotoluene	25.95	8071	101.6000<	100	2%	Acceptable					100	-100%	Fails		(±15)	45
2-Nitrotoluene	22.34	5089	210.9000<	200	5%	Acceptable					200	-100%	Fails		(±15)	45
<b>4-Nitrotoluene</b>				<b>200</b>	<b>-100%</b>	<b>Fails</b>					<b>200</b>	<b>-100%</b>	<b>Fails</b>		(±15)	
3-Nitrotoluene	22.86	3157	100.5000<	100	1%	Acceptable					100	-100%	Fails		(±15)	45
<b>Nitroglycerin</b>				<b>100</b>	<b>-100%</b>	<b>Fails</b>		38.25	5728	101.8000<	100	2%	Acceptable		(±15)	45
<b>PETN</b>				<b>100</b>	<b>-100%</b>	<b>Fails</b>		48.84	10890	105.4000<	100	5%	Acceptable		(±15)	45
3,5-Dinitroaniline	24.62	6884	102.6000<	100	3%	Acceptable					100	-100%	Fails		(±15)	45
<b>EGDN</b>				<b>100</b>	<b>-100%</b>	<b>Fails</b>		17.78	4429	104.7000	100	5%	Acceptable		(±15)	45

Notes M = Manually Integrated Signals Differ by More Than 40%  
 D = Operator Disabled Result Signals Differ by More Than 50%  
 O = Over Calibration Range

2324MW 09MW

A-000014.D

Chromatography Summary

Injection Date: 7/9/2009 3:06 Operator: fhk  
 DataFile: LC10107082009 BVA-000014.D Vial Num: 20  
 Instrument ID: LC10

Method 8330 Target Analyte Results

Sample: LF3H41AC 9187206 G9G030156-1 1X

Method File: LC10107082009 B8330AB.M  
 Start Cal Date: 6/10/2009 15:33 End Cal Date: 6/11/2009 13:47

Matrix: WATER SubList: WATER sub SpikeList:  
 Samp. Info: LF3H41AC 9187206 G9G030156-1 1X,0,  
 Misc. Info: ..1007.3.,20;1;WATER sub.;0;1;LF3H41AC

Dilution Factor	Extract Volume	Sample Volume	Sample Weight
1X	20 mL	1007.3 mL	0 g

Signal 1 UV 250-265						Signal 2 UV 358-205							
Compound Name	RT	Diff	Response	Conc (ug/L)	Flag	RT	Diff	Response	Conc (ug/L)	Flag	MDL	RL	Flag
3,4-Dinitrotoluene	18.24	0.032	5816	2.4480<		18.24	0.032	11243	2.5820		0.0090	0.00	
HMX											0.0268	0.10	
✓ RDX	7.88	0.072	11039	2.8140<	- confirmed on CV						0.0645	0.10	45
Picric ACID											0.1241	0.99	
1,3,5-Trinitrobenzene	10.39	0.189	2271	0.3055<	- Not confirmed on CV						0.0308	0.10	45
1,3-Dinitrobenzene											0.0496	0.10	
TETRYL											0.0496	0.10	
Nitrobenzene											0.0496	0.10	
2,4,6-Trinitrotoluene											0.0238	0.10	
4-AM-2,6-DNT											0.0218	0.10	
2-AM-4,6-DNT											0.0993	0.20	
2,6-Dinitrotoluene											0.0496	0.10	
2,4-Dinitrotoluene											0.0496	0.10	
2-Nitrotoluene											0.0715	0.50	
4-Nitrotoluene											0.0715	0.50	
3 Nitrotoluene											0.0616	0.50	
Nitroglycerin											0.3276	0.65	
PETN											0.2978	0.65	
3,5-Dinitroaniline											0.0248	0.99	

Surrogates:	Spiked	Recovered	% Rec	Spiked	Recovered	% Rec	Limits
3,4-Dinitrotoluene	2.4819	2.4480	99	2.4819	2.5820	104	(79-111)

Notes: M = Manually Integrated Signals Differ by More Than 40%  
 D = Operator Disabled Result Signals Differ by More Than 50%  
 O = Over Calibration Range

Chromatography Summary

Injection Date: 7/11/2009 7:59 Operator: fhk  
 DataFile: LC9.I07102009A.BIC-000015.D Vial Num: 21  
 Instrument ID: LC9

Method 8330 Target Analyte Results

Sample: **LF3H41AC 9187206 G9G030156-1 IX**

Method File: LC9.I07102009A.BIC330CNAB.M  
 Start Cal Date: 6/9/2009 20:53 End Cal Date: 6/10/2009 4:20

Matrix: WATER SubList: WATER sub SpikeList:  
 Samp. Info: LF3H41AC 9187206 G9G030156-1 IX,0,  
 Misc. Info: ..1007 3.,20.1,WATER sub;;0,1;LF3H41AC

Dilution Factor	Extract Volume	Sample Volume	Sample Weight
IX	20 mL	1007.3 mL	0 g

Signal 1 UV 250-265						Signal 2 UV 358-205							
Compound Name	RT	Diff	Response	Conc (ug/L)	Flag	RT	Diff	Response	Conc (ug/L)	Flag	MDI	RL	Flag
3,4-Dinitrotoluene	30.74	0.088	5308	2.6583<							0.0000	0.00	45
HMX											0.0268	0.10	
<input checked="" type="checkbox"/> RDX	25.25	0.116	5549	2.5890<	-confirmed						0.0645	0.10	45
Picric ACID											0.1241	0.99	
1,3,5-Trinitrobenzene											0.0308	0.10	
1,3-Dinitrobenzene											0.0496	0.10	
TETRYL											0.0496	0.10	
Nitrobenzene											0.0496	0.10	
2,4,6-Trinitrotoluene											0.0238	0.10	
4-AM-2,6-DNT	29.18	0.136	110	0.0326<							0.0218	0.10	45
2-AM-4,6-DNT											0.0993	0.20	
2,6-Dinitrotoluene											0.0496	0.10	
2,4-Dinitrotoluene											0.0496	0.10	
2-Nitrotoluene											0.0715	0.49	
4-Nitrotoluene											0.0715	0.49	
3-Nitrotoluene											0.0616	0.49	
Nitroglycerin											0.3276	0.64	
PETN											0.2978	0.64	
3,5-Dinitroaniline											0.0248	0.99	

Surrogates	Spiked	Recovered	% Rec	Spiked	Recovered	% Rec	Limits
3,4-Dinitrotoluene	2.4819	2.6580	107	2.4819		0	(48-143)

Notes: M = Manually Integrated Signals Differ by More Than 40%  
 D = Operator Disabled Result Signals Differ by More Than 50%  
 O = Over Calibration Range

23 LHMW 02 MW

TestAmerica West Sacramento

A-000015.D

Chromatography Summary

Injection Date: 7/9/2009 3:57 Operator: fhk  
Data File: LC10\07082009 BVA-000015.D Vial Num: 21  
Instrument ID: LC10

Method 8330 Target Analyte Results

Sample: LF3H51AC 9187206 G9G030156-2 1X

Method File: LC10\07082009 BVA330AB.M  
Start Cal Date: 6/10/2009 15:33 End Cal Date: 6/11/2009 13:47

Matrix: WATER SubList: WATER sub SpikeList:  
Samp. Info: LF3H51AC 9187206 G9G030156-2 1X,0,  
Misc. Info: :.996 15;;20, L, WATER sub., 0 1:LF3H51AC

Dilution Factor	Extract Volume	Sample Volume	Sample Weight
IX	20 mL	996.15 mL	0 g

Signal 1 UV 250-265

Signal 2 UV 358-205

Compound Name	RT	Diff	Response	Conc (ug/L)	Flag	RT	Diff	Response	Conc (ug/L)	Flag	MDL	RL	Flag
3,4-Dinitrotoluene	18.23	0.022	5996	2.5520 <		18.23	0.022	11570	2.6870		0.0000	0.00	
✓ HMX	5.30	0.005	23018	3.8250 <							0.0271	0.10	45
✓ RDX	7.87	0.018	74461	19.1900 <							0.0653	0.10	45
Picric ACID											0.1255	1.00	
1,3,5-Trinitrobenzene	10.37	0.168	2549	0.3467 <							0.0311	0.10	45
1,3-Dinitrobenzene											0.0502	0.10	
TETRYL											0.0502	0.10	
Nitrobenzene											0.0502	0.10	
2,4,6-Trinitrotoluene											0.0241	0.10	
4-AM-2,6-DNT											0.0221	0.10	
2-AM-4,6-DNT											0.1004	0.20	
2,6-Dinitrotoluene											0.0502	0.10	
2,4-Dinitrotoluene											0.0502	0.10	
2-Nitrotoluene											0.0723	0.50	
4-Nitrotoluene											0.0723	0.50	
3-Nitrotoluene											0.0622	0.50	
Nitroglycerin											0.3313	0.65	
PETN											0.3012	0.65	
3,5-Dinitroaniline											0.0251	1.00	

*Handwritten notes:*  
-> confirmed on cu  
Not confirmed on cu  
7/14/09

Surrogates:	Spiked	Recovered	% Rec	Spiked	Recovered	% Rec	Limits
3,4-Dinitrotoluene	2.5097	2.5520	102	2.5097	2.6870	107	(79-111)

Notes: M = Manually Integrated Signals Differ by More Than 40%  
D = Operator Disabled Result Signals Differ by More Than 50%  
O = Over Calibration Range



Chromatography Summary

Injection Date: 7/11/2009 9:01 Operator: fnk  
 Data File: LC9 (07102009A) BIC-000016.D Vial Num: 22  
 Instrument ID: LC9

Method 8330 Target Analyte Results

Sample: **LF3H51AC 9187206 G9G030156-2 IX**

Method File: LC9 (07102009A) B18330CNAB.M  
 Start Cal Date: 6/9/2009 20:53 End Cal Date: 6/10/2009 4:20

Matrix: WATER SubList: WATER sub Spikelist:  
 Samp. Info: LF3H51AC 9187206 G9G030156-2 IX,0,  
 Misc. Info: .,996.15;;20,1,WATER sub,,0;1;LF3H51AC

Dilution Factor	Extract Volume	Sample Volume	Sample Weight
IX	20 mL	996.15 mL	0 g

Signal 1 UV 250-265						Signal 2 UV 358-205							
Compound Name	RT	Diff	Response	Conc (ug/L)	Flag	RT	Diff	Response	Conc (ug/L)	Flag	MDL	RL	Flag
3,4-Dinitrotoluene	30.72	0.070	5563	2.8170<							0.0000	0.00	45
✓ HMX	34.70	0.046	6845	3.3350<	-> confirmed						0.0271	0.10	45
✓ RDX	25.24	0.104	37128	17.5200<							0.0653	0.10	45
Picric ACID											0.1255	1.01	
1,3,5-Trinitrobenzene											0.0311	0.10	
1,3-Dinitrobenzene											0.0502	0.10	
TETRYL											0.0502	0.10	
Nitrobenzene											0.0502	0.10	
2,4,6-Trinitrotoluene											0.0241	0.10	
1-AM-2,6-DNT	29.10	0.063	107	0.0321<							0.0221	0.10	45
2-AM-4,6-DNT											0.1004	0.20	
2,6-Dinitrotoluene											0.0502	0.10	
2,4-Dinitrotoluene											0.0502	0.10	
2-Nitrotoluene											0.0723	0.50	
4-Nitrotoluene											0.0723	0.50	
3-Nitrotoluene											0.0622	0.50	
Nitroglycerin											0.3313	0.66	
PETN											0.3012	0.66	
3,5-Dinitroaniline											0.0251	1.01	

*Not confirmed ID*

*JKK 7/14/09*

Surrogates:	Spiked	Recovered	% Rec	Spiked	Recovered	% Rec	Limits
3,4-Dinitrotoluene	2.5097	2.8170	112	2.5097	0	0	(48-143)

Notes: M = Manually Integrated Signals Differ by More Than 40%  
 D = Operator Disabled Result Signals Differ by More Than 50%  
 O = Over Calibration Range

2344 MW02BW

TestAmerica West Sacramento

A-000023.D

Chromatography Summary

Injection Date: 7/11/2009 6:11 Operator: fhk  
Data File: LC10 107102009.B\A-000023.D Vial Num: 28  
Instrument ID: LC10

Method 8330 Target Analyte Results

Sample: LF3H61AC 9187206 G9G030156-3 5X

Method File: LC10 107102009 B\8330AB.M  
Start Cal Date: 6/10/2009 15:33 End Cal Date: 6/11/2009 13:47

Matrix: WATER SubList: WATER sub SpikeList:  
Samp. Info: LF3H61AC 9187206 G9G030156-3 5X,0,  
Misc. Info: ;1005.3;20;1;WATER sub;0.5

Dilution Factor	Extract Volume	Sample Volume	Sample Weight
5X	20 mL	1005.3 mL	0 g

Signal 1 UV 250-265						Signal 2 UV 358-205							
Compound Name	RT	Diff	Response	Conc (ug/L)	Flag	RT	Diff	Response	Conc (ug/L)	Flag	MDL	RL	Flag
3,4-Dinitrotoluene	18.34	0.066	1202	2.5350<		18.34	0.069	2416	2.7800		0.0000	0.00	
✓ HMX	5.31	0.016	5180	4.2650<							0.1343	0.50	45
✓ RDX	7.93	0.062	65192	83.2500<							0.3233	0.50	45
Picric ACID											0.6217	4.97	
1,3,5-Trinitrobenzene											0.1542	0.50	
1,3-Dinitrobenzene											0.2487	0.50	
TETRYL											0.2487	0.50	
Nitrobenzene											0.2487	0.50	
2,4,6-Trinitrotoluene											0.1194	0.50	
4-AM-2,6-DNT											0.1094	0.50	
2-AM-4,6-DNT											0.4974	0.99	
2,6-Dinitrotoluene											0.2487	0.50	
✓ 2,4-Dinitrotoluene	21.17	0.042	308	0.3578<							0.2487	0.50	45
2-Nitrotoluene											0.3581	2.49	
4-Nitrotoluene											0.3581	2.49	
3-Nitrotoluene											0.3084	2.49	
Nitroglycerin											1.6413	3.23	
PETN											1.4921	3.23	
3,5-Dinitroaniline											0.1243	4.97	

*continued on cv  
7/14/09*

Surrogates:	Spiked	Recovered	% Rec	Spiked	Recovered	% Rec	Limits
3,4-Dinitrotoluene	2.4868	2.5350	102	2.4868	2.7800	112	(70-111)

Notes: M = Manually Integrated Signals Differ by More Than 40%  
D = Operator Disabled Result Signals Differ by More Than 50%  
O = Over Calibration Range

Chromatography Summary

Injection Date: 7/11/2009 10:03 Operator: fhk  
 Data File: LC9 I07102009A.BIC-000017.D Vial Num: 23  
 Instrument ID: LC9

Method 8330 Target Analyte Results

Sample: **LF3H61AC 9187206 G9G030156-3 5x 5X**

Method File: LC9 I07102009A.BI8330CNAB.M  
 Start Cal Date: 6/9/2009 20:53 End Cal Date: 6/10/2009 4:20

Matrix: WATER SubList: WATER sub SpikeList:  
 Samp. Info: LF3H61AC 9187206 G9G030156-3 5x 5X;0;  
 Misc. Info: ,,1005 3,,20,L,WATER sub;0,5,LF3H61AC

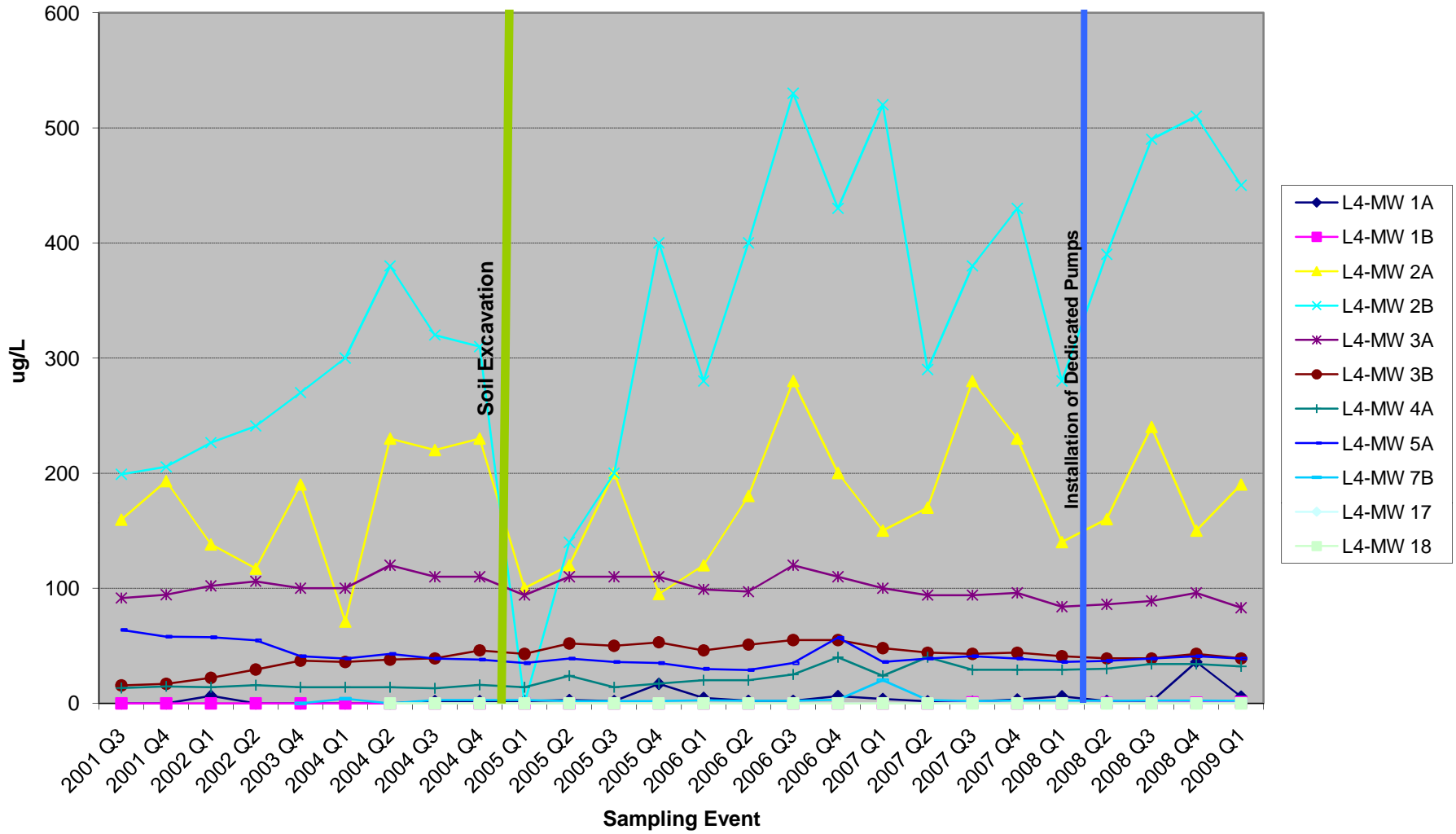
Dilution Factor	Extract Volume	Sample Volume	Sample Weight
5X	20 mL	1005.3 mL	0 g

Signal 1 UV 250-265						Signal 2 UV 358-205							
Compound Name	RT	Diff	Response	Conc (ug/L)	Flag	RT	Diff	Response	Conc (ug/L)	Flag	MDL	RL	Flag
3,4-Dinitrotoluene	30.69	0.036	1177	2.9520<							0.0000	0.00	45
✓ HMX	34.70	0.045	1480	3.5720<							0.1343	0.49	45
✓ RDX	25.20	0.068	31840	74.4200<							0.3233	0.49	45
Picric ACID											0.6217	4.95	
1,3,5-Trinitrobenzene											0.1542	0.49	
1,3-Dinitrobenzene	19.29	0.131	544	0.6145<							0.2487	0.49	45
TETRYL											0.2487	0.49	
Nitrobenzene											0.2487	0.49	
2,4,6-Trinitrotoluene											0.1194	0.49	
4-AM-2,6-DNT											0.1094	0.49	
2-AM-4,6-DNT											0.4974	0.99	
2,6-Dinitrotoluene											0.2487	0.49	
✓ 2,4-Dinitrotoluene	25.95	0.003	348	0.4359<							0.2487	0.49	45
2-Nitrotoluene											0.3581	2.47	
4-Nitrotoluene											0.3581	2.47	
3-Nitrotoluene											0.3084	2.47	
Nitroglycerin											1.6413	3.22	
PETN											1.4921	3.22	
3,5-Dinitroaniline											0.1243	4.95	

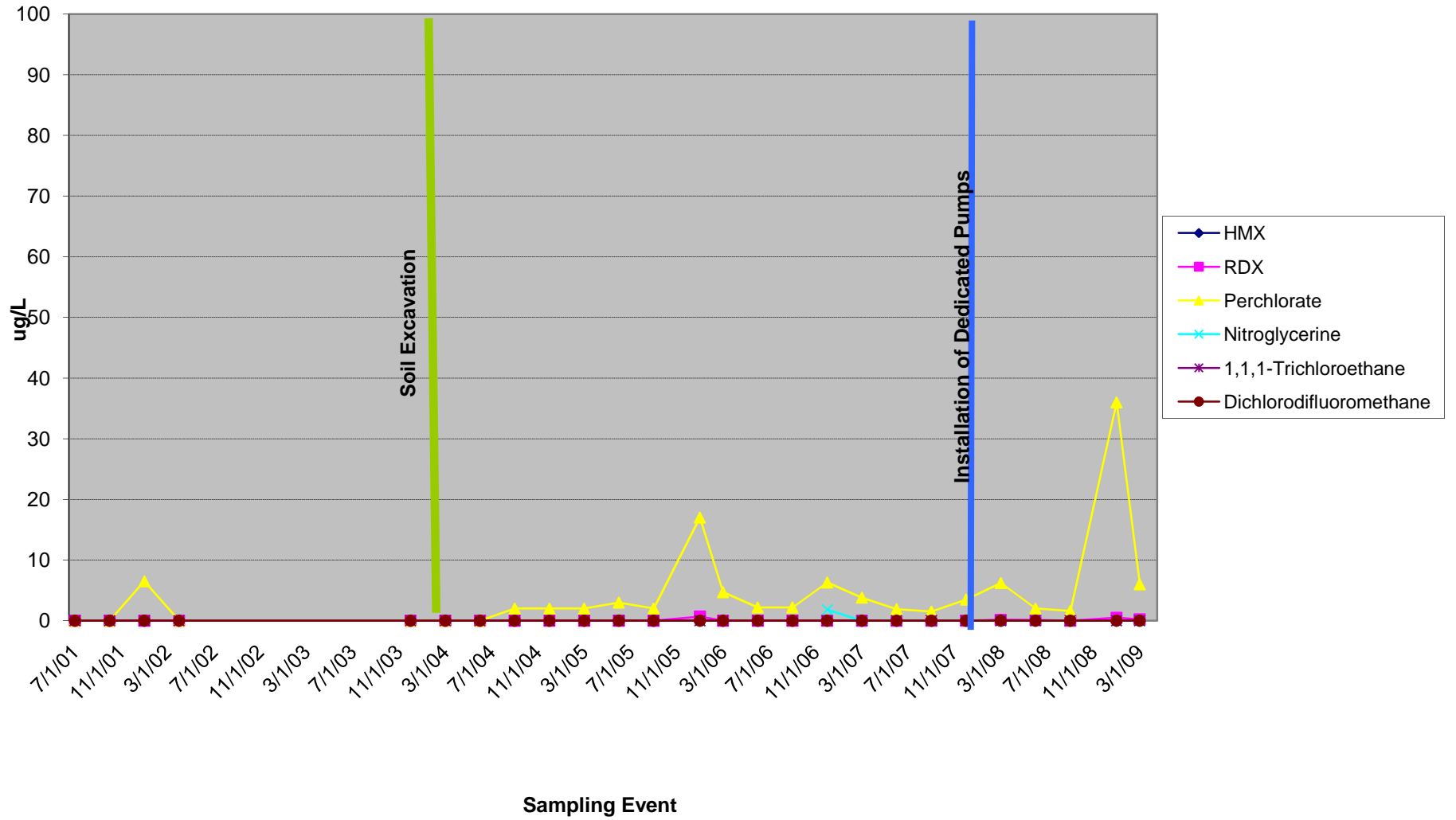
Surrogates:	Spiked	Recovered	% Rec	Spiked	Recovered	% Rec	Limits
3,4-Dinitrotoluene	2.4868	2.9520	119	2.4868	0	0	(48-143)

Notes M = Manually Integrated Signals Differ by More Than 40%  
 D = Operator Disabled Result Signals Differ by More Than 50%  
 O = Over Calibration Range

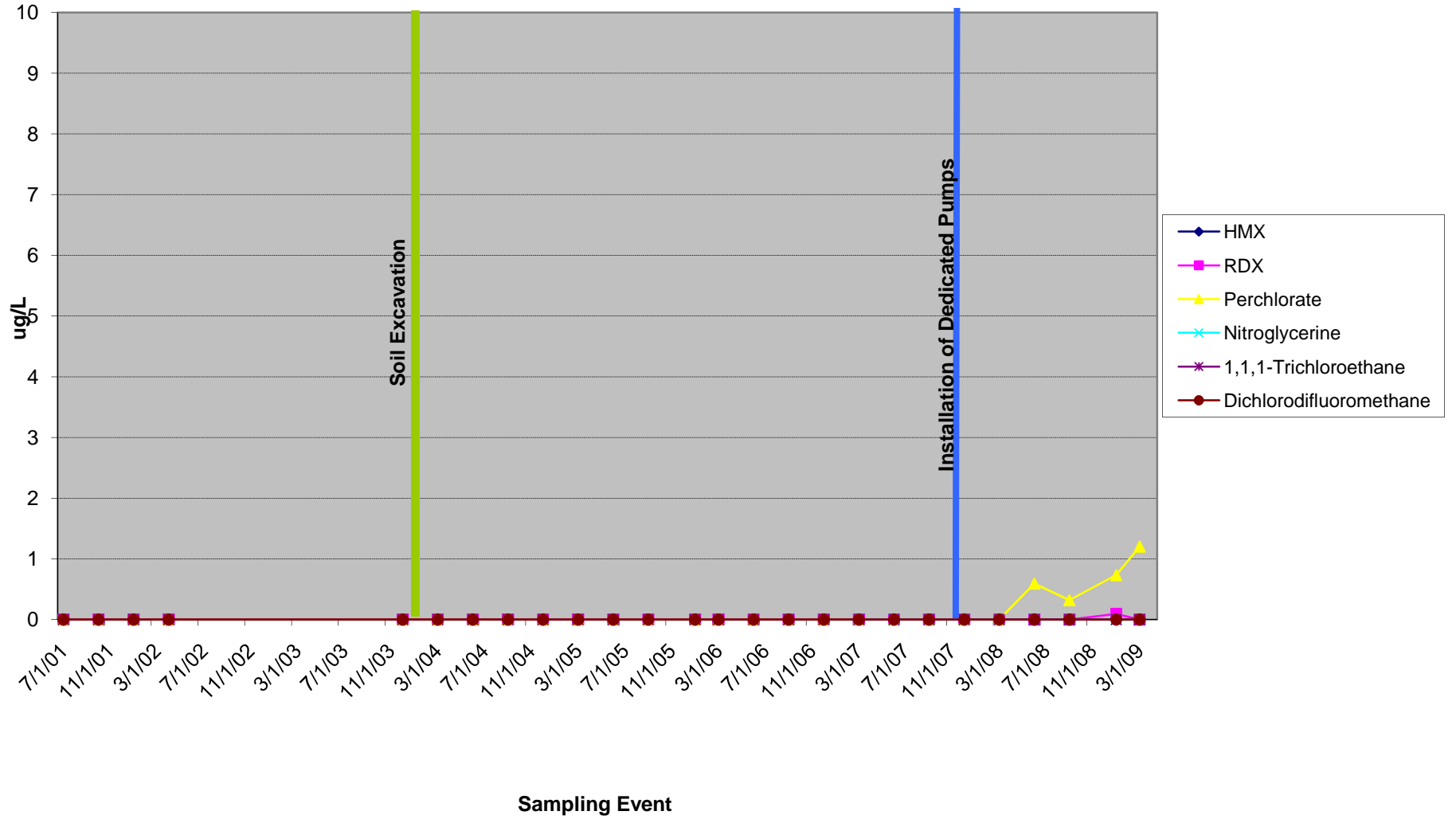
### Landfill 4 Perchlorate Results



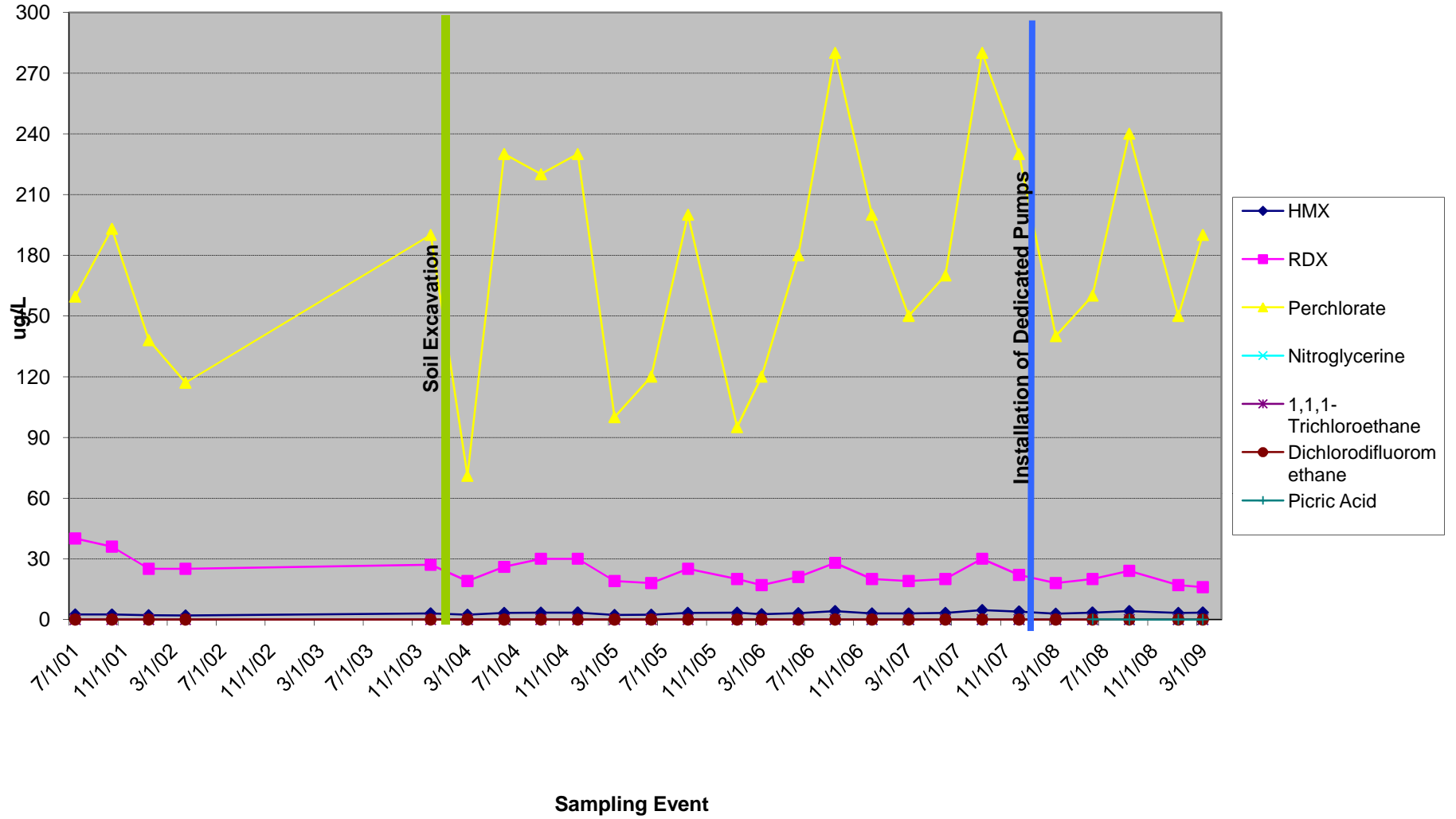
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# L4-MW-1B



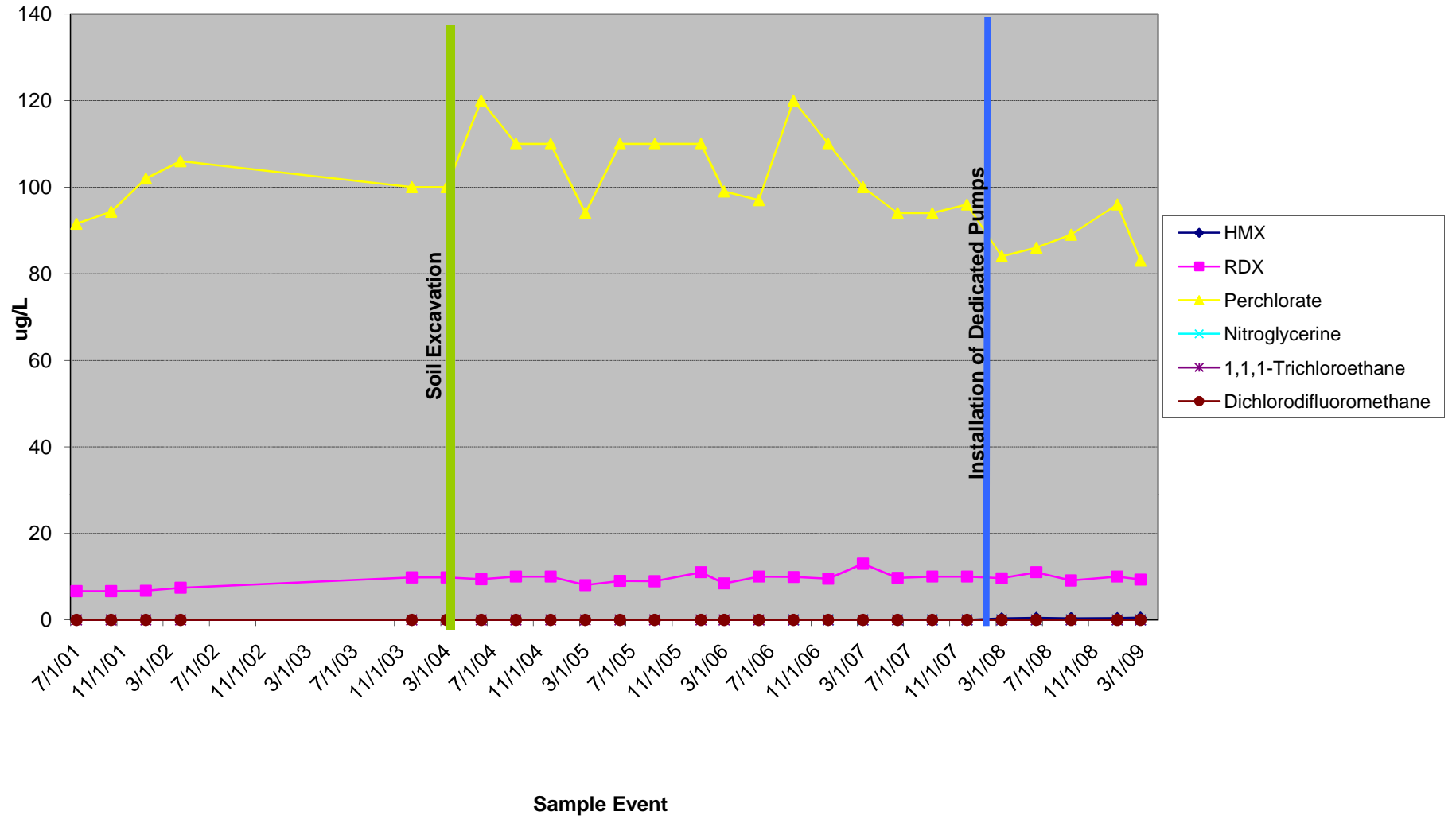
### L4-MW-2A



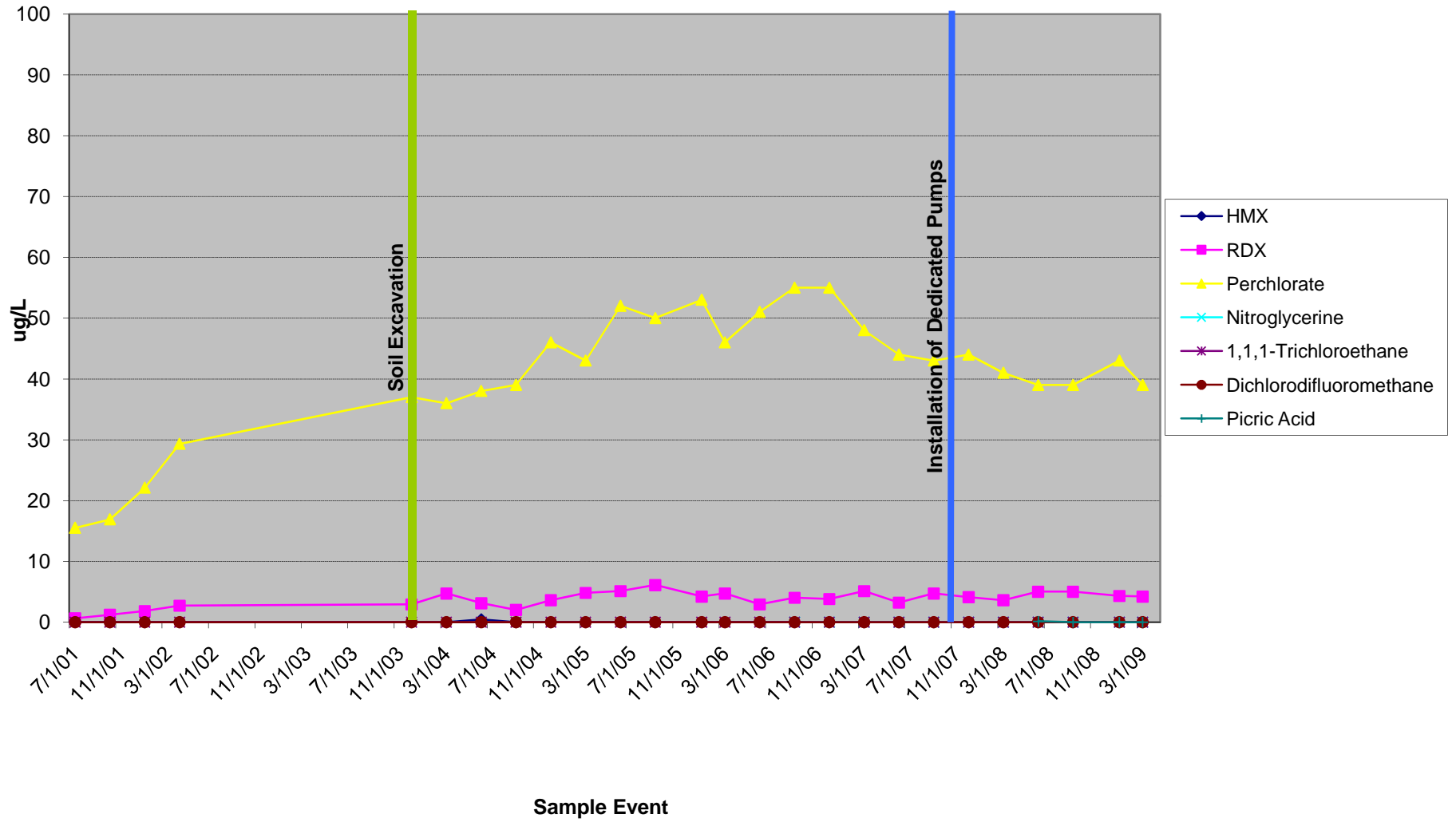




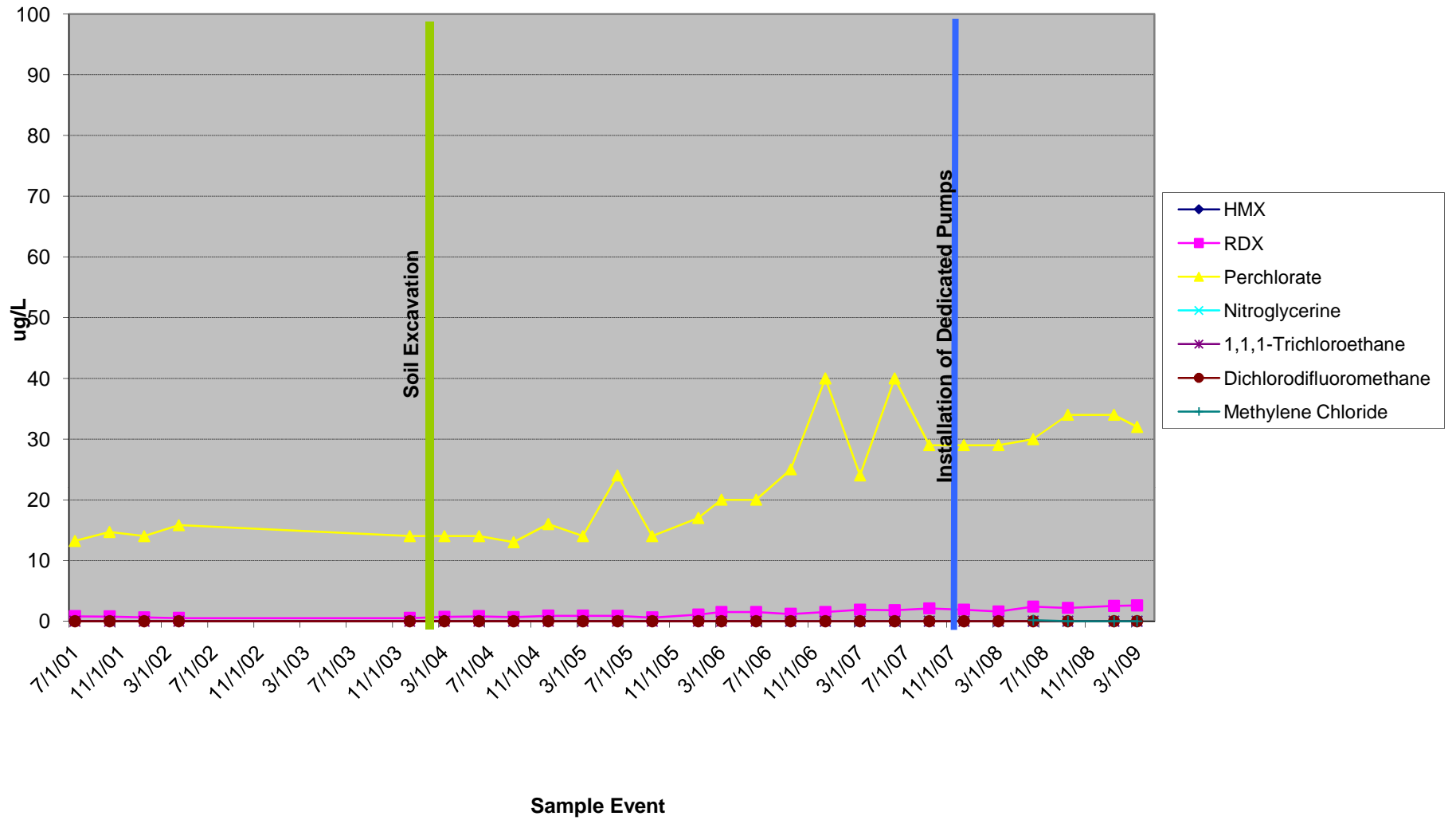
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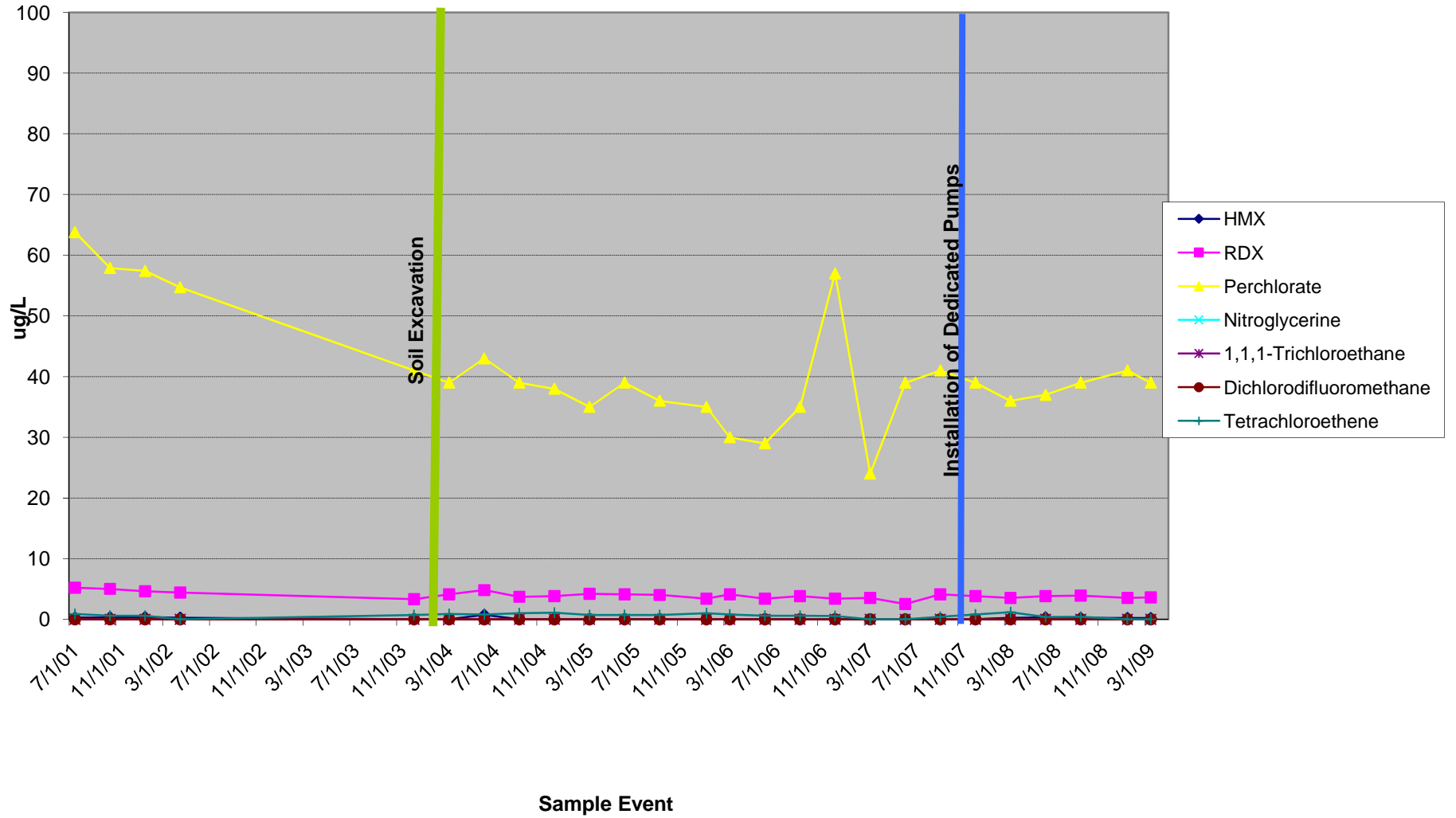
# L4-MW-3B



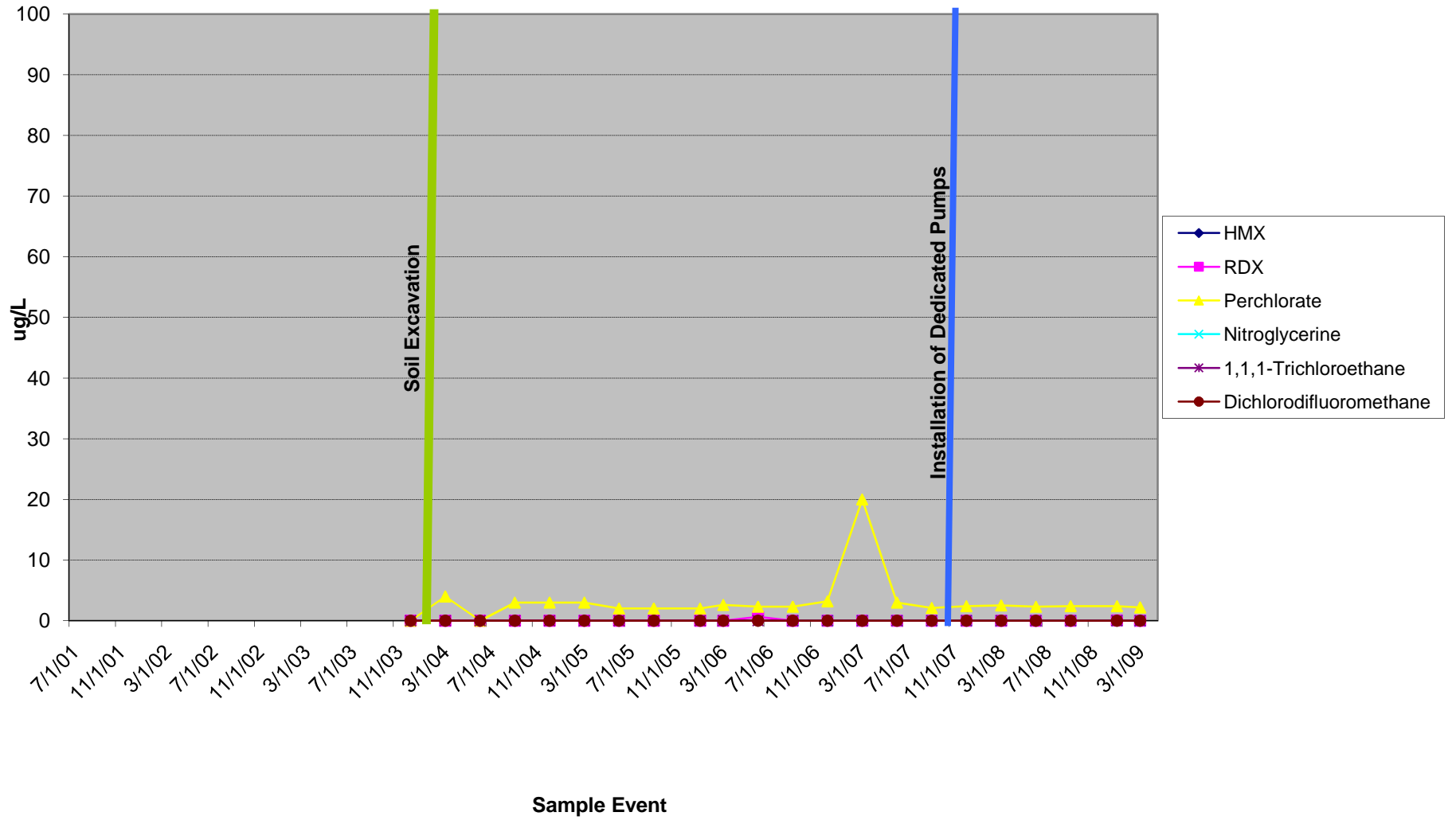
# L4-MW-4A



# L4-MW-5A



# L4-MW-7B





# L4-MW-18

