

DATA VALIDATION SUMMARY REPORT

for Samples Collected During

Groundwater Monitoring

Fort Wingate Depot Activity

McKinley County, New Mexico

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INTRODUCTION

The following data validation summary report covers four (4) water samples, and associated field quality control (QC) samples collected on October 10, 2024, at Fort Wingate Depot Activity (FWDA), located in McKinley County, New Mexico. The samples were logged under Sample Delivery Group (SDG) 280-197922.

The samples in this SDG were analyzed for the following parameters: orthophosphate as P by EPA Method 365.1, anions by U.S. EPA Method 9056A, volatile organic compounds (VOCs) by U.S. EPA Method 8260D, semivolatile organic compounds (SVOCs) by U.S. EPA Method 8270E, total petroleum hydrocarbons-gasoline/diesel/oil range organics (TPH-GRO/DRO/ORO) by U.S. EPA Method 8015D, explosives by EPA Method 8330B, perchlorate by EPA Method 6850, metals (total and dissolved) by U.S. EPA Method 6020B and mercury (total and dissolved) by U.S. EPA Method 7470A.

All samples were collected by Eco & Associates, Inc. (ECO) and were submitted for analysis to Eurofins Environmental Testing America (EETA) Denver located in Arvada, Colorado. All containers were received by EETA at temperatures within the required temperature range of 0.1 to 6.0° Celsius. All containers were received at the laboratory in good condition.

All samples were prepared and analyzed following the procedures outlined in the project-specific Uniform Federal Policy - Quality Assurance Project Plan (UFP-QAPP) and the Department of Defense (DoD) Quality Systems Manual (QSM) Version 5.4. The following table details the samples included in this SDG discussed in this report and the analytical parameters performed.

SAMPLE IDS AND REQUESTED PARAMETERS

| Client Sample ID | Laboratory Sample ID | Matrix | Parameter |
|------------------------------|----------------------|--------|-----------------|
| QC10102024TB (Trip Blank) | 280-197922-1 | Water | V, TPH |
| TMW21102024 | 280-197922-2 | Water | A, O |
| TMW32102024 | 280-197922-3 | Water | A, O |
| TMW38102024 | 280-197922-4 | Water | A, O |
| TMW21102024 | 280-197922-5 | Water | V, M, P, E, TPH |
| TMW32102024 | 280-197922-6 | Water | V, M, P, E |
| TMW38102024 | 280-197922-7 | Water | V, S, M, P, E |

Parameters:

A=Anions

O= Orthophosphate as P

V=VOCs

S=SVOCs

TPH=GRO/DRO/ORO

E=Explosives

P=Perchlorate

M=Metals/Mercury

EXTRACTION, ANALYTICAL, AND REPORTING DETAILS

| Parameter | Matrix | Prep Method | Analytical Method | Units |
|---------------------|--------|-------------|-------------------|-------|
| Anions | Water | -- | SW846 9056A | µg/L |
| Orthophosphate as P | Water | -- | EPA 365.1 | µg/L |
| VOCs | Water | -- | SW846 8260D | µg/L |
| SVOCs | Water | 3510C | SW846 8270E | µg/L |
| TPH GRO | Water | -- | SW846 8015D | µg/L |
| TPH DRO/ORO | Water | 3510C | SW846 8015D | µg/L |
| Explosives | Water | 3535 | EPA 8330B | µg/L |
| Perchlorate | Water | -- | EPA 6850 | µg/L |
| Metals | Water | 3005A/3020A | SW846 6020B | µg/L |
| Mercury | Water | 7470A | SW846 7470A | µg/L |

µg/L= micrograms per liter

EVALUATION CRITERIA

The data submitted by the laboratory has been reviewed and validated at a Stage 2B Validation was performed following the guidelines outlined in the project-specific UFP QAPP, DoD General Data Validation Guidelines, Rev 1 (Nov 2019) and published data validation guideline modules. Information reviewed in the data packages included sample results; field and laboratory quality control results; instrument calibration; calibration verifications; case narratives; sample receipt forms, chain-of-custody (COC) forms. The

analyses and findings presented in this report are based on the reviewed information, and whether guidelines in the associated analytical method, DoD QSM and QAPP were met.

A table detailing the data qualifiers applied for the samples in this SDG as a result of the data validation process is included as Attachment A to this report. Data validation checklists for each analytical method listed in the table above are also included in this report as Attachment B. An ADR.net summary report is included in this report as Attachment C.

ANIONS

General

The anions portion of this SDG consisted of three (3) water samples. The samples were collected on October 10, 2024, and were analyzed for anions as specified in the project-specific UFP-QAPP.

The anions analyses were performed in accordance with U.S. EPA Method SW846 9056A. All samples in this SDG were analyzed following the procedures outlined in the DoD QSM, version 5.4 and the project QAPP.

All samples were analyzed within the holding time required by the method with the following exceptions: sample TMW32102024 was re-analyzed for nitrate outside the 48-hour holding time documented in the QAPP. As such, the nitrate result was qualified “J” as estimated.

Accuracy

Accuracy was evaluated using the percent recovery obtained from the Laboratory Control Sample (LCS), Laboratory Control Sample Duplicate (LCSD), Matrix Spike (MS) and Matrix Spike Duplicate (MSD). Samples TMW21102024 and TMW32102024 were designated for MS/MSD analysis by the laboratory.

All LCS/LCSD and MS/MSD spike recoveries were within acceptance criteria.

Precision

Precision was evaluated using the relative percent difference (RPD) obtained from the LCS/LCSD, MS/MSD and laboratory duplicate concentrations.

All LCS/LCSD and MS/MSD RPDs were within acceptance criteria.

All laboratory duplicate RPDs were within acceptance criteria, except for the following:

| Parent Sample TMW32102024 | | |
|----------------------------------|-------------|-----------------|
| Analyte | %RPD | Criteria |
| Fluoride | 29 | $RPD \leq 10$ |

Fluoride recovered high in the laboratory duplicate. The result in sample TMW32102024 was qualified “J” as estimated.

Representativeness

Representativeness expresses the degree to which sample data accurately and precisely represents actual site conditions. Representativeness has been evaluated by:

- Comparing the COC procedures to those described in the DoD QSM and project QAPP;
- Comparing actual analytical procedures to those described in the DoD QSM and project-specific UFP-QAPP;
- Evaluating holding times; and
- Examining blanks for cross contamination of samples during analysis.

The samples in this SDG were analyzed following the COC and the analytical procedures described in the DoD QSM and project-specific UFP-QAPP. All samples were prepared and analyzed within the holding time required by the method with the exception previously noted. The following QC elements were also evaluated:

- All initial calibration (ICAL) criteria were met.
- The initial calibration verification (ICV) samples were prepared from a second source standard. All ICV criteria were met.
- All initial and continuing calibration blanks (ICB/CCB) criteria were met.
- All continuing calibration verification (CCV) criteria were met.

Six laboratory method blanks were associated with the anions analyses in this SDG. The laboratory method blanks were non-detect for all target anions.

Completeness

Completeness has been evaluated by comparing the total number of samples collected with the total number of samples with valid analytical data.

All results for anions for the samples in this SDG were considered usable. Therefore, the completeness for the anions portion of this SDG is 100%, which meets the minimum acceptance criteria of 90%.

ORTHOPHOSPHATE AS P

General

The orthophosphate portion of this SDG consisted of three (3) water samples. The samples were collected on October 10, 2024 and were analyzed for orthophosphate as specified in the project-specific UFP-QAPP.

The orthophosphate analyses were performed in accordance with U.S. EPA Method 365.1. All samples in this SDG were analyzed following the procedures outlined in the DoD QSM, version 5.4 and the project QAPP. All samples were analyzed within the holding time required by the method.

Accuracy

Accuracy was evaluated using the percent recovery obtained from the LCS/LCSD, and MS/MSD. Sample TMW32102024 was designated for MS/MSD analysis by the laboratory.

All LCS/LCSD and MS/MSD spike recoveries were within acceptance criteria.

Precision

Precision was evaluated using the RPD obtained from the LCS/LCSD and MS/MSD concentrations.

All LCS/LCSD and MS/MSD RPDs were within acceptance criteria.

Representativeness

Representativeness expresses the degree to which sample data accurately and precisely represents actual site conditions. Representativeness has been evaluated by:

- Comparing the COC procedures to those described in the DoD QSM and project QAPP;
- Comparing actual analytical procedures to those described in the DoD QSM and project-specific UFP-QAPP;
- Evaluating holding times; and
- Examining blanks for cross contamination of samples during analysis.

The samples in this SDG were analyzed following the COC and the analytical procedures described in the DoD QSM and project-specific UFP-QAPP. All samples were prepared and analyzed within the holding time required by the method. The following QC elements were also evaluated:

- All ICAL criteria were met.
- The ICV samples were prepared from a second source standard. All ICV criteria were met.
- All ICB/CCB criteria were met.
- All CCV criteria were met.

One laboratory method blank was associated with the orthophosphate analyses in this SDG. The laboratory method blank was non-detect for orthophosphate.

Completeness

Completeness has been evaluated by comparing the total number of samples collected with the total number of samples with valid analytical data.

All results for orthophosphate for the samples in this SDG were considered usable. Therefore, the completeness for the orthophosphate portion of this SDG is 100%, which meets the minimum acceptance criteria of 90%.

VOLATILE ORGANIC COMPOUNDS

General

The VOCs portion of this SDG consisted of four (4) water samples. The samples were collected on October 10, 2024, and were analyzed for VOCs as specified in the project-specific UFP-QAPP.

The VOC analyses were performed in accordance with U.S. EPA Method 8260D. All samples in this SDG were analyzed following the procedures outlined in the DoD QSM, version 5.4 and the project QAPP. All samples were prepared and analyzed within the holding time required by the method. The method requirement for headspace was not met for sample QC10102024TB. As such, all VOCs were qualified “UJ” as estimated at the reporting limit.

Accuracy

Accuracy was evaluated using the percent recovery obtained from the LCS/LCSD and the surrogate spikes. Insufficient sample volume was available to perform a MS/MSD.

All LCS/LCSD recoveries were within acceptance criteria.

Surrogate spike compounds were added to every field and QC sample. All surrogate spike recoveries were within acceptance criteria.

Precision

Precision was evaluated using the RPD obtained from the LCS/LCSD concentrations.

All LCS/LCSD RPDs were within acceptance criteria.

Representativeness

Representativeness expresses the degree to which sample data accurately and precisely represents actual site conditions. Representativeness has been evaluated by:

- Comparing the COC procedures to those described in the DoD QSM and project QAPP;
- Comparing actual analytical procedures to those described in the DoD QSM and project-specific UFP-QAPP;
- Evaluating holding times; and

- Examining blanks for cross contamination of samples during analysis.

The samples in this SDG were analyzed following the COC and the analytical procedures described in the DoD QSM and project-specific UFP-QAPP. All samples were prepared and analyzed within the holding time required by the method. The following QC elements were also evaluated:

- All instrument tune criteria were met.
- All ICAL criteria were met.
- The ICV samples were prepared from a second source standard. All ICV criteria were met.
- All CCV criteria were met.
- All internal standard criteria were met.

Two laboratory method blanks were associated with the VOC analyses in this SDG. The laboratory method blanks were non-detect for all target VOCs.

One trip blank was associated with the VOC analyses in this SDG. The trip blank was non-detect for all target VOCs.

Completeness

Completeness has been evaluated by comparing the total number of samples collected with the total number of samples with valid analytical data.

All results for VOCs for the samples in this SDG were considered usable. Therefore, the completeness for the VOCs portion of this SDG is 100%, which meets the minimum acceptance criteria of 90%.

SEMI-VOLATILE ORGANIC COMPOUNDS

General

The SVOCs portion of this SDG consisted of one (1) water sample. The sample was collected on October 10, 2024, and was analyzed for SVOCs as specified in the project-specific UFP-QAPP.

The SVOC analyses were performed in accordance with U.S. EPA Method 8270E. The sample in this SDG was analyzed following the procedures outlined in the DoD QSM, version 5.4 and the project QAPP. The sample was prepared and analyzed within the holding time required by the method.

Accuracy

Accuracy was evaluated using the percent recovery obtained from the LCS and surrogate spikes.

All LCS spike recoveries were within acceptance criteria.

Surrogate spike compounds were added to every field and QC sample. All surrogate spike recoveries were within acceptance criteria. It should be noted that surrogate 2,4,6-tribromophenol recovery in the CCV associated with batch 671354 was outside control limits. The surrogate recoveries in the associated samples were within control limits; therefore, corrective action was not necessary, and qualification of data was not warranted.

Precision

Precision is evaluated using the RPD obtained from the LCS/LCSD and MS/MSD concentrations. Since a LCSD and MS/MSD were not reported, precision could not be evaluated.

Representativeness

Representativeness expresses the degree to which sample data accurately and precisely represents actual site conditions. Representativeness has been evaluated by:

- Comparing the COC procedures to those described in the DoD QSM and project QAPP;
- Comparing actual analytical procedures to those described in the DoD QSM and project-specific UFP-QAPP;
- Evaluating holding times; and
- Examining blanks for cross contamination of samples during analysis.

The samples in this SDG were analyzed following the COC and the analytical procedures described in the DoD QSM and project-specific UFP-QAPP. The sample was prepared and analyzed within the holding time required by the method. The following QC elements were also evaluated:

- All instrument tune criteria were met.
- All ICAL criteria were met.
- The ICV samples were prepared from a second source standard. All ICV criteria were met.
- All CCV criteria were met
- All internal standard criteria were met. It should be noted that internal standard retention times for the CCV associated with batch 671354 was outside acceptance criteria from the mid-point of the initial calibration. All associated samples were within the acceptance criteria of the daily calibration verification; therefore, corrective action was not necessary, and qualification of data was not warranted.

One laboratory method blank was associated with the SVOC analysis in this SDG. The laboratory method blank was non-detect for all target SVOCs.

Completeness

Completeness has been evaluated by comparing the total number of samples collected with the total number of samples with valid analytical data.

All results for SVOCs for the sample in this SDG were considered usable. Therefore, the completeness for the SVOCs portion of this SDG is 100%, which meets the minimum acceptance criteria of 90%.

TOTAL PETROLEUM HYDROCARBONS GRO

General

The TPH GRO portion of this SDG consisted of two (2) water samples. The samples were collected on October 10, 2024, and were analyzed for TPH GRO as specified in the project-specific UFP-QAPP.

The TPH GRO analyses were performed in accordance with U.S. EPA Method 8015D. All samples in this SDG were analyzed following the procedures outlined in the DoD QSM, version 5.4 and the project QAPP. All samples were prepared and analyzed within the holding time required by the method.

Accuracy

Accuracy was evaluated using the percent recovery obtained from the LCS/LCSD and the surrogate spikes.

All LCS/LCSD spike recoveries were within acceptance criteria.

Surrogate spike compounds were added to every field and QC sample. All surrogate spike recoveries were within acceptance criteria.

Precision

Precision was evaluated using the relative percent difference (RPD) obtained from the LCS/LCSD concentrations.

All LCS/LCSD RPDs were within acceptance criteria.

Representativeness

Representativeness expresses the degree to which sample data accurately and precisely represents actual site conditions. Representativeness has been evaluated by:

- Comparing the COC procedures to those described in the DoD QSM and project QAPP;
- Comparing actual analytical procedures to those described in the DoD QSM and project-specific UFP-QAPP;
- Evaluating holding times; and
- Examining blanks for cross contamination of samples during analysis.

The samples in this SDG were analyzed following the COC and the analytical procedures described in the DoD QSM and project-specific UFP-QAPP. All samples were prepared and analyzed within the holding time required by the method. The following QC elements were also evaluated:

- All ICAL criteria were met.
- The ICV samples were prepared from a second source standard. All ICV criteria were met.
- All CCV criteria were met.

One laboratory method blank was associated with the TPH GRO analyses in this SDG. The laboratory method blank was non-detect for TPH GRO.

One trip blank was associated with the TPH GRO analyses in this SDG. The trip blank was non-detect for TPH-GRO.

Completeness

Completeness has been evaluated by comparing the total number of samples collected with the total number of samples with valid analytical data.

All results for TPH GRO for the samples in this SDG were considered usable. Therefore, the completeness for the TPH GRO portion of this SDG is 100%, which meets the minimum acceptance criteria of 90%.

TOTAL PETROLEUM HYDROCARBONS DRO/ORO

General

The TPH DRO/ORO portion of this SDG consisted of one (1) water sample. The sample was collected on October 10, 2024, and was analyzed for TPH DRO/ORO as specified in the project-specific UFP-QAPP.

The TPH DRO/ORO analysis was performed in accordance with U.S. EPA Method 8015D. The sample in this SDG was analyzed following the procedures outlined in the DoD QSM, version 5.4 and the project QAPP. The sample was prepared and analyzed within the holding time required by the method.

Accuracy

Accuracy was evaluated using the percent recovery obtained from the LCS/LCSD and the surrogate spikes. Insufficient sample volume was available to perform an MS/MSD.

All LCS/LCSD spike recoveries were within acceptance criteria.

Surrogate spike compounds were added to every field and QC sample. All surrogate spike recoveries were within acceptance criteria.

Precision

Precision was evaluated using the RPD obtained from the LCS/LCSD concentrations.

All LCS/LCSD RPDs were within acceptance criteria.

Representativeness

Representativeness expresses the degree to which sample data accurately and precisely represents actual site conditions. Representativeness has been evaluated by:

- Comparing the COC procedures to those described in the DoD QSM and project QAPP;
- Comparing actual analytical procedures to those described in the DoD QSM and project-specific UFP-QAPP;
- Evaluating holding times; and
- Examining blanks for cross contamination of samples during analysis.

The samples in this SDG were analyzed following the COC and the analytical procedures described in the DoD QSM and project-specific UFP-QAPP. All samples were prepared and analyzed within the holding time required by the method. The following QC elements were also evaluated:

- All ICAL criteria were met.
- The ICV samples were prepared from a second source standard. All ICV criteria were met.
- All CCV criteria were met.

One laboratory method blank was associated with the TPH DRO/ORO analysis in this SDG. The laboratory method blank was non-detect for TPH DRO/ORO.

Completeness

Completeness has been evaluated by comparing the total number of samples collected with the total number of samples with valid analytical data.

All results for TPH DRO/ORO for the sample in this SDG are considered usable. Therefore, the completeness for the TPH DRO/ORO portion of this SDG is 100%, which meets the minimum acceptance criteria of 90%.

EXPLOSIVES

General

The explosives portion of this SDG consisted of three (3) water samples. The samples were collected on October 10, 2024, and were analyzed for explosives as specified in the project-specific UFP-QAPP.

The explosives analyses were performed in accordance with U.S. EPA Method 8330B. All samples in this SDG were analyzed following the procedures outlined in the DoD QSM, version 5.4 and the project QAPP. All samples were prepared and analyzed within the holding time required by the method.

Accuracy

Accuracy was evaluated using the percent recovery obtained from the LCS and the surrogate spikes.

All LCS spike recoveries were within acceptance criteria.

Surrogate spike compounds were added to every field and QC sample. All surrogate spike recoveries were within acceptance criteria.

Precision

Precision is evaluated using the RPD obtained from the LCS/LCSD and MS/MSD concentrations. Since a LCSD and MS/MSD were not reported, precision could not be evaluated.

Representativeness

Representativeness expresses the degree to which sample data accurately and precisely represents actual site conditions. Representativeness has been evaluated by:

- Comparing the COC procedures to those described in the DoD QSM and project QAPP;
- Comparing actual analytical procedures to those described in the DoD QSM and project-specific UFP-QAPP;
- Evaluating holding times; and
- Examining blanks for cross contamination of samples during analysis.

The samples in this SDG were analyzed following the COC and the analytical procedures described in the DoD QSM and project-specific UFP-QAPP. All samples were prepared and analyzed within the holding time required by the method. The following QC elements were also evaluated:

- All ICAL criteria were met.
- The ICV samples were prepared from a second source standard. All ICV criteria were met.
- All CCV criteria were met.
- Dual column confirmation for the field samples could not be evaluated because the results were non-detect.

One laboratory method blank was associated with the explosives analyses in this SDG. The laboratory method blank was non-detect for all target explosives.

Completeness

Completeness has been evaluated by comparing the total number of samples collected with the total number of samples with valid analytical data.

All results for explosives for the samples in this SDG were considered usable. Therefore, the completeness for the explosives portion of this SDG is 100%, which meets the minimum acceptance criteria of 90%.

PERCHLORATE

General

The perchlorate portion of this SDG consisted of three (3) water samples. The samples were collected on October 10, 2024, and were analyzed for perchlorate as specified in the project-specific UFP-QAPP.

The perchlorate analyses were performed in accordance with U.S. EPA Method 6850. All samples in this SDG were analyzed following the procedures outlined in the DoD QSM, version 5.4 and the project QAPP. All samples were prepared and analyzed within the holding time required by the method.

Accuracy

Accuracy was evaluated using the percent recovery obtained from the LCS.

All LCS spike recoveries were within acceptance criteria.

Precision

Precision is evaluated using the RPD obtained from the LCS/LCSD and MS/MSD concentrations. Since a LCSD and MS/MSD were not reported, precision could not be evaluated.

Representativeness

Representativeness expresses the degree to which sample data accurately and precisely represents actual site conditions. Representativeness has been evaluated by:

- Comparing the COC procedures to those described in the DoD QSM and project QAPP;
- Comparing actual analytical procedures to those described in the DoD QSM and project-specific UFP-QAPP;
- Evaluating holding times; and
- Examining blanks for cross contamination of samples during analysis.

The samples in this SDG were analyzed following the COC and the analytical procedures described in the DoD QSM and project-specific UFP-QAPP. All samples were prepared and analyzed within the holding time required by the method. The following QC elements were also evaluated:

- All isotope ratio criteria were met.
- All ICAL criteria were met.

- The ICV samples were prepared from a second source standard. All ICV criteria were met.
- All interference check solutions (ICS) were within criteria.
- All CCV criteria were met.
- All initial calibration blank (ICB) criteria were met.
- All continuing calibration blank (CCB) criteria were met.
- All internal standard criteria were met.

One laboratory method blank was associated with the perchlorate analyses in this SDG. The laboratory method blank was non-detect for perchlorate.

Completeness

Completeness has been evaluated by comparing the total number of samples collected with the total number of samples with valid analytical data.

All results for perchlorate for the samples in this SDG were considered usable. Therefore, the completeness for the perchlorate portion of this SDG is 100%, which meets the minimum acceptance criteria of 90%.

METALS

General

The metals portion of this SDG consisted of three (3) water samples. The samples were collected on October 10, 2024, and were analyzed for total and dissolved metals as specified in the project-specific UFP-QAPP.

The metals analyses were performed in accordance with U.S. EPA Method 6020B. All samples in this SDG were analyzed following the procedures outlined in the DoD QSM, version 5.4 and the project QAPP. All samples were prepared and analyzed within the holding time required by the method.

Accuracy

Accuracy was evaluated using the percent recovery obtained from the LCS and MS/MSD. Sample TMW38102024 was designated for MS/MSD analysis by the laboratory.

All LCS spike recoveries were within acceptance criteria.

All MS/MSD recoveries were within acceptance criteria. It should be noted that one or more MS/MSD RECs for calcium and sodium exceeded acceptance criteria, however; the sample concentrations were greater than 4 times the MS/MSD spike concentrations. As such, the MS/MSD RECs could not be evaluated, and qualification was not warranted.

Precision

Precision was evaluated using the RPD obtained from the MS/MSD concentrations.

All MS/MSD RPDs were within acceptance criteria.

Representativeness

Representativeness expresses the degree to which sample data accurately and precisely represents actual site conditions. Representativeness has been evaluated by:

- Comparing the COC procedures to those described in the DoD QSM and project QAPP;
- Comparing actual analytical procedures to those described in the DoD QSM and project-specific UFP-QAPP;
- Evaluating holding times; and
- Examining blanks for cross contamination of samples during analysis.

The samples in this SDG were analyzed following the COC and the analytical procedures described in the DoD QSM and project-specific UFP-QAPP. All samples were prepared and analyzed within the holding time required by the method. The following QC elements were also evaluated:

- All instrument tune criteria were met.
- All initial calibration criteria were met.
- The ICV was prepared from a second source standard. All ICV criteria were met.
- All CCV criteria were met.
- All low-level ICV (LL ICV) criteria were met.
- All ICS were within criteria.
- All ICB criteria were met except for the following:
 - The ICB associated with batch 670767 had detections of silver and thallium. Thallium was non-detect in the associated samples, as such, no qualification was warranted. Silver was detected less than 5 times the ICB detection in associated sample TMW32102024, as such, the result was qualified “U” as non-detect.
- All CCB criteria were met except for the following:
 - The CCBs associated with batch 670767 had detections of silver, thallium and magnesium. Thallium and magnesium were either non-detect or greater than 5 times the CCB detections. As such, no qualification was warranted. Silver was detected less than 5 times the CCB detection in associated sample TMW32102024, as such, the results were qualified “U” as non-detect.
- All internal standard criteria associated with the target metals were met.
- A serial dilution test (DT) was performed on the same sample as the MS/MSD. The DT was only applicable for those metals that failed in the MS/MSD and were

detected in the parent sample at a concentration of 50 times the LOQ or greater. All applicable metals met criteria in the DT.

- The post digestion spike (PDS) was performed on the same sample as the MS/MSD. The PDS was only applicable for those metals that failed in the MS/MSD. All PDS recoveries were within acceptance criteria. It should be noted that the PDS REC for calcium exceeded acceptance criteria, however; the sample concentration was greater than 4 times the PDS spike concentration. As such, the PDS REC could not be evaluated, and qualification was not warranted.

Three laboratory method blanks were associated with the metals analyses in this SDG. Aluminum, iron and magnesium were detected in one or more of the laboratory method blanks. The associated samples with results less than 5 times the method blank detections were qualified “U” as non-detect.

Completeness

Completeness has been evaluated by comparing the total number of samples collected with the total number of samples with valid analytical data.

All results for metals for the samples in this SDG were considered usable. Therefore, the completeness for the metals portion of this SDG is 100%, which meets the minimum acceptance criteria of 90%.

MERCURY

General

The mercury portion of this SDG consisted of three (3) water samples. The samples were collected on October 10, 2024, and were analyzed for total and dissolved mercury as specified in the project-specific UFP-QAPP.

The mercury analyses were performed in accordance with U.S. EPA Method 7470A. All samples in this SDG were analyzed following the procedures outlined in the DoD QSM, version 5.4 and the project QAPP. All samples were prepared and analyzed within the holding time required by the method.

Accuracy

Accuracy was evaluated using the percent recovery obtained from the LCS/LCSD.

All LCS/LCSD spike recoveries were within acceptance criteria.

Precision

Precision was evaluated using the RPD obtained from the LCS/LCSD concentrations.

All LCS/LCSD RPDs were within acceptance criteria.

Representativeness

Representativeness expresses the degree to which sample data accurately and precisely represents actual site conditions. Representativeness has been evaluated by:

- Comparing the COC procedures to those described in the DoD QSM and project QAPP;
- Comparing actual analytical procedures to those described in the DoD QSM and project-specific UFP-QAPP;
- Evaluating holding times; and
- Examining blanks for cross contamination of samples during analysis.

The samples in this SDG were analyzed following the COC and the analytical procedures described in the DoD QSM and project-specific UFP-QAPP. All samples were prepared and analyzed within the holding time required by the method. The following QC elements were also evaluated:

- All initial calibration criteria were met.
- The ICV was prepared from a second source standard. All ICV criteria were met.
- All CCV criteria were met.
- All LL ICV criteria were met.
- All ICB criteria were met.
- All CCB criteria were met.

Two laboratory method blanks were associated with the mercury analyses in this SDG. The laboratory method blanks were non-detect for mercury.

Completeness

Completeness has been evaluated by comparing the total number of samples collected with the total number of samples with valid analytical data.

All results for mercury for the samples in this SDG were considered usable. Therefore, the completeness for the mercury portion of this SDG is 100%, which meets the minimum acceptance criteria of 90%.

COMPARABILITY

All data was generated using contract-specific standard methods and reported with known data quality, type of analysis, units, etc.

DATA USABILITY

The purpose of this data validation report is to ensure the integrity and reliability of analytical laboratory data. The data quality is evaluated based on precision, accuracy, representativeness, comparability, and completeness (PARCC) characteristics of the data.

The validated data indicated that the laboratory correctly performed the analyses. Based on the data quality assessment, none of the data were qualified as rejected.

All data in this SDG are considered usable, as qualified, for the purposes of this project. All Method Quality Objectives have been met.

SENSITIVITY

The detection limit (DL), LOD and limit of quantitation (LOQ) values reported for the samples were compared to those listed in WS #15, Table 15.1 of the QAPP to ensure that sensitivity requirements were met. The DL, LOD, and LOQ values matched those listed in the QAPP before dilutions were taken into account. The following LOQs exceed the project quantitation limits (PQLs):

| Methods | Analyte(s) | Sample(s) |
|---------|--|-----------------------------|
| 8270E | 2,4-dinitrotoluene, 2,6-dinitrotoluene, 3,3'-dichlorobenzidine, 4,6-dinitro-2-methylphenol, 4-chloroaniline, benzo[a]anthracene, benzo[a]pyrene, benzo[b]fluoranthene, bis(2-chloroethyl)ether, bis(2-ethylhexyl) phthalate, dibenz(a,h)anthracene, hexachlorobenzene, hexachlorobutadiene, hexachlorocyclopentadiene, hexachloroethane, indeno[1,2,3-cd]pyrene, nitrobenzene, n-nitrosodi-n-propylamine, pentachlorophenol and phenol | TMW38102024 |
| 8015D | DRO | TMW21102024 |
| 8330B | nitroglycerin | TMW21102024 and TMW38102024 |

DATA QUALIFIER CHANGES

The sample results and final data qualifiers and reason codes that were added, removed, or changed as a result of the data validation process are included in a table as Attachment A to this report.

DATA QUALIFIER DEFINITIONS

The data qualifiers are defined in WS #36, Table 36.2 of the project QAPP as follows.

U = The analyte was not detected and was reported as less than the LOD. The LOD has been adjusted for any dilution or concentration of the sample.

J = The reported result was an estimated value with an unknown bias.

J+ = The reported result was an estimated quantity, but the result may be biased high.

J- = The reported result was an estimated quantity, but the result may be biased low.

UJ = The analyte was not detected and was reported as less than the LOD. However, the reported numerical value is approximate.

X= The sample results (including non-detects) were affected by serious deficiencies in the ability to analyze the sample and to meet published method and project quality control criteria. The presence or absence of the analyte cannot be substantiated by the data provided. Acceptance (J-flag) or rejection (R-flag) of the data should be decided by the project team (which should include a project chemist) during the Data Usability Assessment process.

REASON CODE DEFINITIONS

These data validation reason codes were used to document the logic behind all data validation qualifiers:

| Validation Qualifier Reason Codes | Validation Comments |
|-----------------------------------|--|
| BLL | Concentration in equipment blank at or above reporting limit. |
| BLM | Concentration in equipment blank less than reporting limit |
| BLN | Concentration in trip blank at or above reporting limit. |
| BLO | Concentration in trip blank less than reporting limit. |
| BLR | Concentration in field blank or decon blank at or above reporting limit. |
| BLS | Concentration in field blank or decon blank less than reporting limit. |
| BLT | Concentration in method blank less than reporting limit. |
| BLU | Concentration in method blank at or above reporting limit. |
| CO1 | Column confirmation RPD exceeds acceptance limit. |
| CR1 | Result exceeded calibration range. |
| DU1 | Field duplicate RPD exceeds acceptance limit. |
| DU2 | Laboratory duplicate RPD exceeds acceptance limit. |
| DU3 | Field Duplicate RPD not calculated but results demonstrate a high degree of variability. |
| HS | VOA vial has headspace greater than 6 millimeters. |
| LC1 | LCS and/or LCSD recovery above upper acceptance limit. |
| LC2 | LCS and/or LCSD recovery below lower acceptance limit. |
| LC7 | LCS/LCSD RPD exceeds acceptance limit. |
| MD1 | MS and/or MSD recovery above upper acceptance limit. |
| MD2 | MS and/or MSD recovery below lower acceptance limit. |
| MD5 | MS/MSD RPD exceeds acceptance limit. |
| PJ | Professional judgment used. See specific details in Data Validation Report. |
| SC1 | Analysis holding time exceeded. |
| SC3 | Extraction holding time exceeded. |
| SC6 | Temperature of sample outside acceptance range. |
| SU1 | Surrogate recovery above upper acceptance limit. |
| SU2 | Surrogate recovery below lower acceptance limit. |
| PR1 | Samples not properly preserved. |
| TR | Result is detected between the reporting limit and detection limit. |

ACRONYMS AND ABBREVIATIONS

The following is a list of acronyms and abbreviations that were used in this data validation report.

| | |
|----------|---|
| CCB | Continuing Calibration Blank |
| CCV | Continuing Calibration Verification |
| CoC | Chain of Custody |
| DL | Detection Limit |
| DoD | Department of Defense |
| DT | Dilution Test |
| ETTA | Eurofins Environment Testing America |
| FWDA | Fort Wingate Depot Activity |
| ICAL | Initial Calibration |
| ICB | Initial Calibration Blank |
| ICS | Interference Check Sample |
| ICV | Initial Calibration Verification |
| LCS | Laboratory Control Sample |
| LCSD | Laboratory Control Sample Duplicate |
| LOD | Limit of Detection |
| LOQ | Limit of Quantitation |
| MS | Matrix Spike |
| MSD | Matrix Spike Duplicate |
| PQL | Project Quantitation Limit |
| PDS | Post Digestion Spike |
| QC | Quality Control |
| QSM | Quality Systems Manual |
| RPD | Relative Percent Difference |
| SDG | Sample Delivery Group |
| UFP-QAPP | Uniform Federal Policy – Quality Assurance Project Plan |

Attachment A

Validated Data Summary

**Fort Wingate Depot Activity Northern Area
McKinley County, New Mexico
Northern Area Groundwater Sampling
Validated Data Summary for Water Samples Collected September and October 2024**

| SAMPLE ID: | | PROJECT QUANTITATION LIMIT GOAL (PQLG) ^[1] | TMW21102024 | | TMW32102024 | | TMW38102024 | |
|-----------------------------|------|--|------------------------------|---|------------------------------|---|------------------------------|---|
| DATE SAMPLED: | | | 10/10/2024 | | 10/10/2024 | | 10/10/2024 | |
| LAB SAMPLE ID: | | | 280-197922-5 280-197922-2 | | 280-197922-6 280-197922-3 | | 280-197922-7 280-197922-4 | |
| Volatile Organics - SW8260D | | Unit | | | | | | |
| 1,1,1,2-Tetrachloroethane | µg/L | 5.7 | 1.0 | U | 1.0 | U | 1.0 | U |
| 1,1,1-Trichloroethane | µg/L | 200 | 1.0 | U | 1.0 | U | 1.0 | U |
| 1,1,2,2-Tetrachloroethane | µg/L | 10 | 1.0 | U | 1.0 | U | 1.0 | U |
| 1,1,2-Trichloroethane | µg/L | 5 | 1.0 | U | 1.0 | U | 1.0 | U |
| 1,1-Dichloroethane | µg/L | 25 | 1.0 | U | 1.0 | U | 1.0 | U |
| 1,1-Dichloroethene | µg/L | 7 | 1.0 | U | 1.0 | U | 1.0 | U |
| 1,1-Dichloropropene | µg/L | 4.7 | 1.0 | U | 1.0 | U | 1.0 | U |
| 1,2,3-Trichlorobenzene | µg/L | 7 | 4.0 | U | 4.0 | U | 4.0 | U |
| 1,2,3-Trichloropropane | µg/L | 2.5 | 2.5 | U | 2.5 | U | 2.5 | U |
| 1,2,4-Trichlorobenzene | µg/L | 70 | 1.0 | U | 1.0 | U | 1.0 | U |
| 1,2,4-Trimethylbenzene | µg/L | 56 | 1.0 | U | 1.0 | U | 1.0 | U |
| 1,2-Dibromo-3-chloropropane | µg/L | 5 | 5.0 | U | 5.0 | U | 5.0 | U |
| 1,2-Dibromoethane (EDB) | µg/L | 1 | 1.0 | U | 1.0 | U | 1.0 | U |
| 1,2-Dichlorobenzene | µg/L | 600 | 1.0 | U | 1.0 | U | 1.0 | U |
| 1,2-Dichloroethane | µg/L | 5 | 1.0 | U | 1.0 | U | 1.0 | U |
| 1,2-Dichloropropane | µg/L | 5 | 1.0 | U | 1.0 | U | 1.0 | U |
| 1,3,5-Trimethylbenzene | µg/L | 60 | 1.0 | U | 1.0 | U | 1.0 | U |
| 1,3-Dichlorobenzene | µg/L | 75 | 1.0 | U | 1.0 | U | 1.0 | U |
| 1,3-Dichloropropane | µg/L | 370 | 1.0 | U | 1.0 | U | 1.0 | U |
| 1,4-Dichlorobenzene | µg/L | 75 | 1.0 | U | 1.0 | U | 1.0 | U |
| 2,2-Dichloropropane | µg/L | 5 | 1.0 | U | 1.0 | U | 1.0 | U |
| 2-Butanone (MEK) | µg/L | 5,600 | 10 | U | 10 | U | 10 | U |
| 2-Chlorotoluene | µg/L | 240 | 1.0 | U | 1.0 | U | 1.0 | U |
| 2-Hexanone | µg/L | 38 | 5.0 | U | 5.0 | U | 5.0 | U |
| 4-Chlorotoluene | µg/L | 250 | 1.0 | U | 1.0 | U | 1.0 | U |
| 4-Isopropyltoluene | µg/L | 450 | 1.0 | U | 1.0 | U | 1.0 | U |
| 4-Methyl-2-pentanone (MIBK) | µg/L | 6,300 | 5.0 | U | 5.0 | U | 5.0 | U |
| Acetone | µg/L | 18,000 | 15 | U | 15 | U | 15 | U |
| Benzene | µg/L | 5 | 1.0 | U | 1.0 | U | 1.0 | U |
| Bromobenzene | µg/L | 62 | 1.0 | U | 1.0 | U | 1.0 | U |
| Bromochloromethane | µg/L | 83 | 1.0 | U | 1.0 | U | 1.0 | U |
| Bromodichloromethane | µg/L | 80 | 1.0 | U | 1.0 | U | 1.0 | U |
| Bromoform | µg/L | 80 | 2.0 | U | 2.0 | U | 2.0 | U |
| Bromomethane | µg/L | 7.5 | 5.0 | U | 5.0 | U | 5.0 | U |

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Northern Area Groundwater Sampling
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|--------------------------------|------|--|------------------------------|---|------------------------------|---|------------------------------|---|
| DATE SAMPLED: | | | 10/10/2024 | | 10/10/2024 | | 10/10/2024 | |
| LAB SAMPLE ID: | | | 280-197922-5 280-197922-2 | | 280-197922-6 280-197922-3 | | 280-197922-7 280-197922-4 | |
| Carbon disulfide | µg/L | 810 | 2.0 | U | 2.0 | U | 2.0 | U |
| Carbon tetrachloride | µg/L | 5 | 1.0 | U | 1.0 | U | 1.0 | U |
| Chlorobenzene | µg/L | 100 | 1.0 | U | 1.0 | U | 1.0 | U |
| Chloroethane | µg/L | 8,300 | 2.0 | U | 2.0 | U | 2.0 | U |
| Chloroform | µg/L | 80 | 1.0 | U | 1.0 | U | 1.0 | U |
| Chloromethane | µg/L | 190 | 2.0 | U | 2.0 | U | 2.0 | U |
| cis-1,2-Dichloroethene | µg/L | 70 | 1.0 | U | 1.0 | U | 1.0 | U |
| cis-1,3-Dichloropropene | µg/L | 4.7 | 1.0 | U | 1.0 | U | 1.0 | U |
| Dibromochloromethane | µg/L | 80 | 1.0 | U | 1.0 | U | 1.0 | U |
| Dibromomethane | µg/L | 8.3 | 1.0 | U | 1.0 | U | 1.0 | U |
| Dichlorodifluoromethane | µg/L | 200 | 2.0 | U | 2.0 | U | 2.0 | U |
| Ethylbenzene | µg/L | 700 | 1.0 | U | 1.0 | U | 1.0 | U |
| Hexachlorobutadiene | µg/L | 2 | 2.0 | U | 2.0 | U | 2.0 | U |
| Isopropylbenzene | µg/L | 450 | 1.0 | U | 1.0 | U | 1.0 | U |
| Methyl acetate | µg/L | 20,000 | 5.0 | U | 5.0 | U | 5.0 | U |
| Methyl tert-butyl ether (MTBE) | µg/L | 100 | 5.0 | U | 5.0 | U | 5.0 | U |
| Methylene chloride | µg/L | 5 | 2.0 | U | 2.0 | U | 2.0 | U |
| m-Xylene & p-Xylene | µg/L | 620 | 2.0 | U | 2.0 | U | 2.0 | U |
| Naphthalene | µg/L | 30 | 3.0 | U | 3.0 | U | 3.0 | U |
| n-Butylbenzene | µg/L | 1,000 | 1.0 | U | 1.0 | U | 1.0 | U |
| n-Propylbenzene | µg/L | 660 | 1.0 | U | 1.0 | U | 1.0 | U |
| o-Xylene | µg/L | 620 | 1.0 | U | 1.0 | U | 1.0 | U |
| sec-Butylbenzene | µg/L | 2,000 | 1.0 | U | 1.0 | U | 1.0 | U |
| Styrene | µg/L | 100 | 1.0 | U | 1.0 | U | 1.0 | U |
| tert-Butylbenzene | µg/L | 690 | 1.0 | U | 1.0 | U | 1.0 | U |
| Tetrachloroethene | µg/L | 5 | 1.0 | U | 1.0 | U | 1.0 | U |
| Toluene | µg/L | 1,000 | 1.0 | U | 1.0 | U | 1.0 | U |
| trans-1,2-Dichloroethene | µg/L | 100 | 1.0 | U | 1.0 | U | 1.0 | U |
| trans-1,3-Dichloropropene | µg/L | 4.7 | 1.0 | U | 1.0 | U | 1.0 | U |
| Trichloroethene | µg/L | 5 | 1.0 | U | 1.0 | U | 1.0 | U |
| Trichlorofluoromethane | µg/L | 5,200 | 2.0 | U | 2.0 | U | 2.0 | U |
| Vinyl chloride | µg/L | 2 | 1.0 | U | 1.0 | U | 1.0 | U |

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|---------------------------------|------|--|------------------------------|------------------------------|------------------------------|
| DATE SAMPLED: | | | 10/10/2024 | 10/10/2024 | 10/10/2024 |
| LAB SAMPLE ID: | | | 280-197922-5 280-197922-2 | 280-197922-6 280-197922-3 | 280-197922-7 280-197922-4 |
| Semivolatile Organics - SW8270E | | | | | |
| 2,2'-Oxybis (1-chloropropane) | µg/L | 710 | -- | -- | 11 U |
| 2,4,5-Trichlorophenol | µg/L | 1,200 | -- | -- | 11 U |
| 2,4,6-Trichlorophenol | µg/L | 12 | -- | -- | 11 U |
| 2,4-Dichlorophenol | µg/L | 46 | -- | -- | 11 U |
| 2,4-Dimethylphenol | µg/L | 360 | -- | -- | 11 U |
| 2,4-Dinitrophenol | µg/L | 39 | -- | -- | 34 U |
| 2,4-Dinitrotoluene | µg/L | 10 | -- | -- | 11 U |
| 2,6-Dinitrotoluene | µg/L | 10 | -- | -- | 11 U |
| 2-Chloronaphthalene | µg/L | 750 | -- | -- | 4.6 U |
| 2-Chlorophenol | µg/L | 91 | -- | -- | 11 U |
| 2-Methylnaphthalene | µg/L | 30 | -- | -- | 4.6 U |
| 2-Methylphenol | µg/L | 930 | -- | -- | 11 U |
| 2-Nitroaniline | µg/L | 190 | -- | -- | 11 U |
| 2-Nitrophenol | µg/L | na | -- | -- | 11 U |
| 3 & 4 Methylphenol | µg/L | 370 | -- | -- | 11 U |
| 3,3'-Dichlorobenzidine | µg/L | 50 | -- | -- | 57 U |
| 3-Nitroaniline | µg/L | 38 | -- | -- | 11 U |
| 4,6-Dinitro-2-methylphenol | µg/L | 50 | -- | -- | 57 U |
| 4-Bromophenyl phenyl ether | µg/L | na | -- | -- | 11 U |
| 4-Chloro-3-methylphenol | µg/L | 1,400 | -- | -- | 11 U |
| 4-Chloroaniline | µg/L | 20 | -- | -- | 23 U |
| 4-Chlorophenyl phenyl ether | µg/L | na | -- | -- | 11 U |
| 4-Nitroaniline | µg/L | 38 | -- | -- | 11 U |
| 4-Nitrophenol | µg/L | na | -- | -- | 29 U |
| Acenaphthene | µg/L | 530 | -- | -- | 4.6 U |
| Acenaphthylene | µg/L | 120 | -- | -- | 4.6 U |
| Anthracene | µg/L | 1,800 | -- | -- | 4.6 U |
| Benzaldehyde | µg/L | 190 | -- | -- | 5.7 U |
| Benz(a)anthracene | µg/L | 4 | -- | -- | 4.6 U |
| Benzo(a)pyrene | µg/L | 4 | -- | -- | 4.6 U |
| Benzo(b)fluoranthene | µg/L | 4 | -- | -- | 4.6 U |
| Benzo(g,h,i)perylene | µg/L | 120 | -- | -- | 4.6 U |

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|--------------------------------------|------|--|------------------------------|------------------------------|------------------------------|
| DATE SAMPLED: | | | 10/10/2024 | 10/10/2024 | 10/10/2024 |
| LAB SAMPLE ID: | | | 280-197922-5 280-197922-2 | 280-197922-6 280-197922-3 | 280-197922-7 280-197922-4 |
| Benzo(k)fluoranthene | µg/L | 25 | -- | -- | 4.6 U |
| bis(2-Chloroethoxy)methane | µg/L | 59 | -- | -- | 11 U |
| bis(2-Chloroethyl)ether | µg/L | 10 | -- | -- | 11 U |
| bis(2-Ethylhexyl)phthalate | µg/L | 10 | -- | -- | 11 U |
| Butyl benzyl phthalate | µg/L | 160 | -- | -- | 4.6 U |
| Caprolactam | µg/L | 9,900 | -- | -- | 17 U |
| Carbazole | µg/L | 290 | -- | -- | 4.6 U |
| Chrysene | µg/L | 250 | -- | -- | 4.6 U |
| Dibenz(a,h)anthracene | µg/L | 10 | -- | -- | 11 U |
| Dibenzofuran | µg/L | 7.9 | -- | -- | 4.6 U |
| Diethyl phthalate | µg/L | 15,000 | -- | -- | 4.6 U |
| Dimethyl phthalate | µg/L | na | -- | -- | 4.6 U |
| Di-n-butyl phthalate | µg/L | 900 | -- | -- | 4.6 U |
| Di-n-octyl phthalate | µg/L | 200 | -- | -- | 11 U |
| Fluoranthene | µg/L | 800 | -- | -- | 4.6 U |
| Fluorene | µg/L | 290 | -- | -- | 4.6 U |
| Hexachlorobenzene | µg/L | 10 | -- | -- | 11 U |
| Hexachlorobutadiene | µg/L | 10 | -- | -- | 11 U |
| Hexachlorocyclopentadiene | µg/L | 50 | -- | -- | 57 U |
| Hexachloroethane | µg/L | 10 | -- | -- | 11 U |
| Indeno(1,2,3-cd)pyrene | µg/L | 10 | -- | -- | 11 U |
| Isophorone | µg/L | 780 | -- | -- | 11 U |
| Naphthalene | µg/L | 30 | -- | -- | 4.6 U |
| Nitrobenzene | µg/L | 10 | -- | -- | 11 U |
| n-Nitrosodi-n-propylamine | µg/L | 10 | -- | -- | 11 U |
| N-Nitrosodiphenylamine | µg/L | 120 | -- | -- | 11 U |
| Pentachlorophenol | µg/L | 50 | -- | -- | 57 U |
| Phenanthrene | µg/L | 170 | -- | -- | 4.6 U |
| Phenol | µg/L | 10 | -- | -- | 11 U |
| Pyrene | µg/L | 120 | -- | -- | 11 U |
| Petroleum Hydrocarbons - SW8015D | | | | | |
| Gasoline Range Organics (GRO) C6-C10 | µg/L | 25 | 25 U | -- | -- |
| Diesel Range Organics (DRO) C10-C28 | µg/L | 250 | 260 U | -- | -- |
| Oil Range Organics (ORO) C20-C38 | µg/L | 60,200 | 510 U | -- | -- |

**Fort Wingate Depot Activity Northern Area
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|--|------|--|------------------------------|---|------------------------------|---|------------------------------|---|
| DATE SAMPLED: | | | 10/10/2024 | | 10/10/2024 | | 10/10/2024 | |
| LAB SAMPLE ID: | | | 280-197922-5 280-197922-2 | | 280-197922-6 280-197922-3 | | 280-197922-7 280-197922-4 | |
| Explosives - SW8330B | | | | | | | | |
| 1,3,5-Trinitrobenzene | µg/L | 590 | 0.22 | U | 0.21 | U | 0.24 | U |
| 1,3-Dinitrobenzene | µg/L | 2 | 0.11 | U | 0.11 | U | 0.12 | U |
| 2,4,6-Trinitrotoluene (TNT) | µg/L | 9.8 | 0.11 | U | 0.11 | U | 0.12 | U |
| 2,4-Dinitrotoluene | µg/L | 2.4 | 0.10 | U | 0.10 | U | 0.11 | U |
| 2,6-Dinitrotoluene | µg/L | 0.49 | 0.10 | U | 0.10 | U | 0.11 | U |
| 2-Amino-4,6-dinitrotoluene | µg/L | 1.9 | 0.11 | U | 0.11 | U | 0.12 | U |
| 4-Amino-2,6-dinitrotoluene | µg/L | 1.9 | 0.15 | U | 0.15 | U | 0.17 | U |
| Hexahydro-1,3,5-trinitro-1,3,5-triazine (RDX) | µg/L | 9.7 | 0.22 | U | 0.21 | U | 0.24 | U |
| m-Nitrotoluene | µg/L | 1.7 | 0.41 | U | 0.41 | U | 0.45 | U |
| Nitrobenzene | µg/L | 1.4 | 0.22 | U | 0.21 | U | 0.24 | U |
| Nitroglycerin | µg/L | 2.1 | 2.2 | U | 2.1 | U | 2.4 | U |
| Octahydro-1,3,5,7-tetranitro-1,3,5,7-tetrazocine (HMX) | µg/L | 1,000 | 0.22 | U | 0.21 | U | 0.24 | U |
| o-Nitrotoluene | µg/L | 3.1 | 0.22 | U | 0.21 | U | 0.24 | U |
| Pentaerythritol Tetranitrate (PETN) | µg/L | 170 | 1.1 | U | 1.1 | U | 1.2 | U |
| p-Nitrotoluene | µg/L | 43 | 0.42 | U | 0.42 | U | 0.46 | U |
| Trinitrophenylmethylnitramine (Tetryl) | µg/L | 39 | 0.11 | U | 0.11 | U | 0.12 | U |
| Perchlorate - SW6850 | | | | | | | | |
| Perchlorate | µg/L | 14 | 0.20 | U | 420 | | 0.20 | U |
| Metals, Total - SW6020B/SW7470A | | | | | | | | |
| Aluminum | µg/L | 200 | 46 | J | 25 | J | 18 | J |
| Antimony | µg/L | 6 | 2.0 | U | 0.45 | J | 2.0 | U |
| Arsenic | µg/L | 10 | 0.93 | J | 3.5 | J | 0.85 | J |
| Barium | µg/L | 2,000 | 17 | | 7.3 | | 13 | |
| Beryllium | µg/L | 4 | 1.0 | U | 1.0 | U | 1.0 | U |
| Cadmium | µg/L | 5 | 1.0 | U | 0.22 | J | 1.0 | U |
| Calcium | µg/L | na | 34,000 | | 13,000 | | 23,000 | |
| Chromium | µg/L | 50 | 1.3 | J | 0.65 | J | 0.61 | J |
| Cobalt | µg/L | 50 | 1.0 | U | 1.0 | U | 1.0 | U |
| Copper | µg/L | 1,000 | 1.5 | J | 0.86 | J | 1.1 | J |
| Iron | µg/L | 300 | 200 | U | 200 | U | 200 | U |
| Lead | µg/L | 15 | 1.0 | U | 1.0 | U | 1.0 | U |
| Magnesium | µg/L | na | 6,400 | | 1,300 | | 2,500 | |
| Manganese | µg/L | 50 | 0.71 | J | 27 | | 90 | |
| Mercury | µg/L | 2 | 0.20 | U | 0.20 | U | 0.20 | U |
| Nickel | µg/L | 200 | 2.5 | J | 3.0 | U | 3.0 | U |
| Potassium | µg/L | na | 430 | J | 760 | J | 1,000 | |
| Selenium | µg/L | 50 | 8.3 | | 2.8 | J | 5.0 | U |
| Silver | µg/L | 50 | 1.0 | U | 1.0 | U | 1.0 | U |
| Sodium | µg/L | na | 670,000 | | 770,000 | | 960,000 | |
| Thallium | µg/L | 2 | 1.0 | U | 1.0 | U | 1.0 | U |
| Vanadium | µg/L | 86 | 3.2 | J | 2.1 | J | 5.0 | U |
| Zinc | µg/L | 5,000 | 10 | U | 4.8 | J | 2.0 | J |

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|-------------------------------------|------|--|------------------------------|------------------------------|------------------------------|
| DATE SAMPLED: | | | 10/10/2024 | 10/10/2024 | 10/10/2024 |
| LAB SAMPLE ID: | | | 280-197922-5 280-197922-2 | 280-197922-6 280-197922-3 | 280-197922-7 280-197922-4 |
| Metals, Dissolved - SW6020B/SW7470A | | | | | |
| Aluminum | µg/L | 200 | 200 U | 200 U | 200 U |
| Antimony | µg/L | 6 | 2.0 U | 0.40 J | 2.0 U |
| Arsenic | µg/L | 10 | 0.95 J | 1.4 J | 0.66 J |
| Barium | µg/L | 2,000 | 17 | 6.3 | 12 |
| Beryllium | µg/L | 4 | 1.0 U | 1.0 U | 1.0 U |
| Cadmium | µg/L | 5 | 1.0 U | 1.0 U | 1.0 U |
| Calcium | µg/L | na | 32,000 | 13,000 | 21,000 |
| Chromium | µg/L | 50 | 1.3 J | 3.0 U | 3.0 U |
| Cobalt | µg/L | 50 | 1.0 U | 1.0 U | 1.0 U |
| Copper | µg/L | 1,000 | 1.4 J | 2.0 U | 2.0 U |
| Iron | µg/L | 300 | 19 J | 8.8 J | 79 J |
| Lead | µg/L | 15 | 1.0 U | 1.0 U | 1.0 U |
| Magnesium | µg/L | na | 6,600 | 1,300 | 2,300 |
| Manganese | µg/L | 50 | 3.0 U | 28 | 77 |
| Mercury | µg/L | 2 | 0.20 U | 0.20 U | 0.20 U |
| Nickel | µg/L | 200 | 3.0 U | 3.0 U | 3.0 U |
| Potassium | µg/L | na | 450 J | 860 J | 1,100 |
| Selenium | µg/L | 50 | 8.0 | 3.1 J | 5.0 U |
| Silver | µg/L | 50 | 1.0 U | 1.0 U | 1.0 U |
| Sodium | µg/L | na | 690,000 | 820,000 | 990,000 |
| Thallium | µg/L | 2 | 1.0 U | 1.0 U | 1.0 U |
| Vanadium | µg/L | 86 | 2.9 J | 2.2 J | 5.0 U |
| Zinc | µg/L | 5,000 | 4.4 J | 10 U | 10 |

**Fort Wingate Depot Activity Northern Area
McKinley County, New Mexico
Northern Area Groundwater Sampling
Validated Data Summary for Water Samples Collected September and October 2024**

| SAMPLE ID: | | PROJECT QUANTITATION LIMIT GOAL (PQLG) ^[1] | TMW21102024 | TMW32102024 | TMW38102024 |
|---------------------------------|------|--|------------------------------|------------------------------|------------------------------|
| DATE SAMPLED: | | | 10/10/2024 | 10/10/2024 | 10/10/2024 |
| LAB SAMPLE ID: | | | 280-197922-5 280-197922-2 | 280-197922-6 280-197922-3 | 280-197922-7 280-197922-4 |
| General Chemistry | | | | | |
| Orthophosphate as P - EPA 365.1 | | | | | |
| Orthophosphate as P | µg/L | 20,000 | 38 J | 50 U | 39 J |
| Anions - SW9056A | | | | | |
| Bromide | µg/L | na | 600 | 850 | 620 |
| Chloride | µg/L | 250,000 | 71,000 | 180,000 | 480,000 |
| Fluoride | µg/L | 1,600 | 1,000 | 940 J | 1,600 |
| Nitrate as N | µg/L | 10,000 | 12,000 | 3,000 J- | 500 U |
| Nitrite as N | µg/L | 1,000 | 500 U | 500 U | 500 U |
| Sulfate | µg/L | 250,000 | 540,000 | 1,100,000 | 1,300,000 |

QA NOTES AND DATA QUALIFIERS:

* - Field duplicate of sample on left.
(NO CODE) - Confirmed identification.
U - Analyte was analyzed for but not detected above the reported limit of quantitation (LOQ).
UJ - Analyte not detected, reported LOQ may be inaccurate or imprecise.
J - Analyte detected, estimated concentration.
J- - Analyte detected, estimated concentration with a low bias.
J+ - Analyte detected, estimated concentration with a high bias.
X - The presence or absence of the analyte cannot be substantiated due to deficiencies in meeting QC criteria.
Detections are bolded.

Detections above the PQLG are highlighted.

NOTES:

[1] The PQLG is the lower of the New Mexico Water Quality Control Commission standard (NM WQCC) and the EPA MCL. If the analyte does not have an NM WQCC or MCL but has an EPA Tap Water RSL, the lower value between the adjusted carcinogenic RSL (target excess cancer risk level of 1 x 10⁻⁵) and the non-carcinogenic RSL (with a target hazard index of 1.0) was selected.

µg/L - micrograms per liter

na - Limit not available

-- Analyte was not tested.

Fort Wingate Depot Activity Northern Area
McKinley County, New Mexico
Northern Area Groundwater Sampling
Validated Data Summary for Water QC Samples Collected September
and October 2024

| SAMPLE ID: | | QC10102024TB | |
|------------------------------------|------|--------------|----|
| DATE SAMPLED: | | 10/10/2024 | |
| LAB SAMPLE ID: | | 280-197922-1 | |
| | Unit | | |
| Volatile Organics - SW8260D | | | |
| 1,1,1,2-Tetrachloroethane | µg/L | 1.0 | UJ |
| 1,1,1-Trichloroethane | µg/L | 1.0 | UJ |
| 1,1,2,2-Tetrachloroethane | µg/L | 1.0 | UJ |
| 1,1,2-Trichloroethane | µg/L | 1.0 | UJ |
| 1,1-Dichloroethane | µg/L | 1.0 | UJ |
| 1,1-Dichloroethene | µg/L | 1.0 | UJ |
| 1,1-Dichloropropene | µg/L | 1.0 | UJ |
| 1,2,3-Trichlorobenzene | µg/L | 4.0 | UJ |
| 1,2,3-Trichloropropane | µg/L | 2.5 | UJ |
| 1,2,4-Trichlorobenzene | µg/L | 1.0 | UJ |
| 1,2,4-Trimethylbenzene | µg/L | 1.0 | UJ |
| 1,2-Dibromo-3-chloropropane | µg/L | 5.0 | UJ |
| 1,2-Dibromoethane (EDB) | µg/L | 1.0 | UJ |
| 1,2-Dichlorobenzene | µg/L | 1.0 | UJ |
| 1,2-Dichloroethane | µg/L | 1.0 | UJ |
| 1,2-Dichloropropane | µg/L | 1.0 | UJ |
| 1,3,5-Trimethylbenzene | µg/L | 1.0 | UJ |
| 1,3-Dichlorobenzene | µg/L | 1.0 | UJ |
| 1,3-Dichloropropane | µg/L | 1.0 | UJ |
| 1,4-Dichlorobenzene | µg/L | 1.0 | UJ |
| 2,2-Dichloropropane | µg/L | 1.0 | UJ |
| 2-Butanone (MEK) | µg/L | 10 | UJ |
| 2-Chlorotoluene | µg/L | 1.0 | UJ |
| 2-Hexanone | µg/L | 5.0 | UJ |
| 4-Chlorotoluene | µg/L | 1.0 | UJ |
| 4-Isopropyltoluene | µg/L | 1.0 | UJ |
| 4-Methyl-2-pentanone (MIBK) | µg/L | 5.0 | UJ |
| Acetone | µg/L | 15 | UJ |
| Benzene | µg/L | 1.0 | UJ |
| Bromobenzene | µg/L | 1.0 | UJ |
| Bromochloromethane | µg/L | 1.0 | UJ |
| Bromodichloromethane | µg/L | 1.0 | UJ |
| Bromoform | µg/L | 2.0 | UJ |
| Bromomethane | µg/L | 5.0 | UJ |
| Carbon disulfide | µg/L | 2.0 | UJ |
| Carbon tetrachloride | µg/L | 1.0 | UJ |
| Chlorobenzene | µg/L | 1.0 | UJ |
| Chloroethane | µg/L | 2.0 | UJ |
| Chloroform | µg/L | 1.0 | UJ |
| Chloromethane | µg/L | 2.0 | UJ |
| cis-1,2-Dichloroethene | µg/L | 1.0 | UJ |
| cis-1,3-Dichloropropene | µg/L | 1.0 | UJ |
| Dibromochloromethane | µg/L | 1.0 | UJ |
| Dibromomethane | µg/L | 1.0 | UJ |
| Dichlorodifluoromethane | µg/L | 2.0 | UJ |
| Ethylbenzene | µg/L | 1.0 | UJ |
| Hexachlorobutadiene | µg/L | 2.0 | UJ |
| Isopropylbenzene | µg/L | 1.0 | UJ |
| Methyl acetate | µg/L | 5.0 | UJ |
| Methyl tert-butyl ether (MTBE) | µg/L | 5.0 | UJ |

Fort Wingate Depot Activity Northern Area
McKinley County, New Mexico
Northern Area Groundwater Sampling
Validated Data Summary for Water QC Samples Collected September
and October 2024

| SAMPLE ID: | | QC10102024TB | |
|---|------|--------------|----|
| DATE SAMPLED: | | 10/10/2024 | |
| LAB SAMPLE ID: | | 280-197922-1 | |
| Methylene chloride | µg/L | 2.0 | UJ |
| m-Xylene & p-Xylene | µg/L | 2.0 | UJ |
| Naphthalene | µg/L | 3.0 | UJ |
| n-Butylbenzene | µg/L | 1.0 | UJ |
| n-Propylbenzene | µg/L | 1.0 | UJ |
| o-Xylene | µg/L | 1.0 | UJ |
| sec-Butylbenzene | µg/L | 1.0 | UJ |
| Styrene | µg/L | 1.0 | UJ |
| tert-Butylbenzene | µg/L | 1.0 | UJ |
| Tetrachloroethene | µg/L | 1.0 | UJ |
| Toluene | µg/L | 1.0 | UJ |
| trans-1,2-Dichloroethene | µg/L | 1.0 | UJ |
| trans-1,3-Dichloropropene | µg/L | 1.0 | UJ |
| Trichloroethene | µg/L | 1.0 | UJ |
| Trichlorofluoromethane | µg/L | 2.0 | UJ |
| Vinyl chloride | µg/L | 1.0 | UJ |
| Petroleum Hydrocarbons - SW8015D | | | |
| Gasoline Range Organics (GRO) C6-C10 | µg/L | 25 | U |
| Diesel Range Organics (DRO) C10-C28 | µg/L | -- | |
| Oil Range Organics (ORO) C20-C38 | µg/L | -- | |

QA NOTES AND DATA QUALIFIERS:

(NO CODE) - Confirmed identification.

U - Analyte was analyzed for but not detected above the reported limit of detection (LOQ).

UJ - Analyte not detected, reported LOQ may be inaccurate or imprecise.

J - Analyte detected, estimated concentration.

Detections are bolded.

NOTES:

µg/L - micrograms per liter

-- Analyte was not tested.

Attachment B

Checklists

VALIDATION CHECKLIST

SDG#: 280-197922

Date: 2/14/25

Laboratory: EETA

Reviewer: Kortney Curry

Method: 6020 & 7470

| | Validation Area | Acceptable? Y/N/NA | Comments |
|-------|--|-----------------------|----------|
| I. | Case narrative | Y | |
| II. | Sample receipt/Technical holding times | Y | |
| III. | Instrument performance check/Tune | Y | |
| IV. | Initial calibration/ICV/LLICV | Y | |
| V. | Continuing Calibration | Y | |
| VI. | Laboratory Blanks- MB, ICB/CCB | N | See DVR |
| VI. | Field blanks | NA | |
| VII. | Interference check standard | Y | |
| VIII. | Matrix spike/Matrix spike duplicate | Y | |
| IX. | Laboratory control samples | Y | |
| X. | Field duplicates/Field triplicates | NA | |
| XI. | Internal standards | Y | |
| XII. | Dilution test | Y | |
| XIII. | Post digestion spike | Y | |
| XIV. | Compound quantitation LOQ/LOD/DL | Y | |
| XV. | Target compound identification | Y | |

VALIDATION CHECKLIST

SDG#: 280-197922

Date: 2/14/25

Laboratory: EETA

Reviewer: Kortney Curry

Method: 6850

| | Validation Area | Acceptable? Y/N/NA | Comments |
|-------|--|-----------------------|----------|
| I. | Case narrative | Y | |
| II. | Sample receipt/Technical holding times | Y | |
| III. | Instrument performance check/Tune | NA | |
| IV. | Initial calibration/ICV/LLICV | Y | |
| V. | Continuing Calibration | Y | |
| VI. | Laboratory Blanks- MB, ICB/CCB | Y | |
| VI. | Field blanks | NA | |
| VII. | Surrogate spikes | NA | |
| VIII. | Matrix spike/Matrix spike duplicate | NA | |
| IX. | Laboratory control samples | Y | |
| X. | Field duplicates/Field triplicates | NA | |
| XI. | Internal standards | Y | |
| XII. | Compound quantitation LOQ/LOD/DL | Y | |
| XIII. | Target compound identification | Y | |

VALIDATION CHECKLIST

SDG#: 280-197922

Date: 2/14/25

Laboratory: EETA

Reviewer: Kortney Curry

Method: 8260

| | Validation Area | Acceptable? Y/N/NA | Comments |
|-------|--|-----------------------|----------|
| I. | Case narrative | Y | |
| II. | Sample receipt/Technical holding times | N | See DVR |
| III. | Instrument performance check/Tune | Y | |
| IV. | Initial calibration/ICV | Y | |
| V. | Continuing Calibration | Y | |
| VI. | Laboratory Blanks- MB | Y | |
| VI. | Field blanks | Y | |
| VII. | Surrogate spikes | Y | |
| VIII. | Matrix spike/Matrix spike duplicate | NA | |
| IX. | Laboratory control samples | Y | |
| X. | Field duplicates/Field triplicates | NA | |
| XI. | Internal standards | Y | |
| XII. | Compound quantitation LOQ/LOD/DL | Y | |
| XIII. | Target compound identification | Y | |

VALIDATION CHECKLIST

SDG#: 280-197922

Date: 2/14/25

Laboratory: EETA

Reviewer: Kortney Curry

Method: 8270

| | Validation Area | Acceptable? Y/N/NA | Comments |
|-------|--|-----------------------|----------|
| I. | Case narrative | Y | |
| II. | Sample receipt/Technical holding times | Y | |
| III. | Instrument performance check/Tune | Y | |
| IV. | Initial calibration/ICV | Y | |
| V. | Continuing Calibration | Y | |
| VI. | Laboratory Blanks- MB | Y | |
| VI. | Field blanks | Y | |
| VII. | Surrogate spikes | Y | |
| VIII. | Matrix spike/Matrix spike duplicate | NA | |
| IX. | Laboratory control samples | Y | |
| X. | Field duplicates/Field triplicates | NA | |
| XI. | Internal standards | Y | |
| XII. | Compound quantitation LOQ/LOD/DL | Y | |
| XIII. | Target compound identification | Y | |

VALIDATION CHECKLIST

SDG#: 280-197922

Date: 2/14/25

Laboratory: EETA

Reviewer: Kortney Curry

Method: 8330

| | Validation Area | Acceptable? Y/N/NA | Comments |
|-------|--|-----------------------|----------|
| I. | Case narrative | Y | |
| II. | Sample receipt/Technical holding times | Y | |
| III. | Instrument performance check/Tune | NA | |
| IV. | Initial calibration/ICV | Y | |
| V. | Continuing Calibration | Y | |
| VI. | Laboratory Blanks- MB, ICB/CCB | Y | |
| VI. | Field blanks | NA | |
| VII. | Interference check standard | NA | |
| VIII. | Matrix spike/Matrix spike duplicate | NA | |
| IX. | Laboratory control samples | Y | |
| X. | Lab duplicates | NA | |
| XI. | External standards | Y | |
| XII. | Dilution test | NA | |
| XIII. | Post digestion spike | NA | |
| XIV. | Compound quantitation LOQ/LOD/DL | Y | |
| XV. | Target compound identification | Y | |

VALIDATION CHECKLIST

SDG#: 280-197922

Date: 2/14/25

Laboratory: EETA

Reviewer: Kortney Curry

Method: 9056A & 365.1

| | Validation Area | Acceptable? Y/N/NA | Comments |
|-------|--|-----------------------|----------|
| I. | Case narrative | Y | |
| II. | Sample receipt/Technical holding times | N | See DVR |
| III. | Instrument performance check/Tune | NA | |
| IV. | Initial calibration/ICV | Y | |
| V. | Continuing Calibration | Y | |
| VI. | Laboratory Blanks- MB, ICB/CCB | Y | |
| VI. | Field blanks | NA | |
| VII. | Interference check standard | NA | |
| VIII. | Matrix spike/Matrix spike duplicate | Y | |
| IX. | Laboratory control samples | Y | |
| X. | Lab duplicates | N | See DVR |
| XI. | External standards | NA | |
| XII. | Dilution test | NA | |
| XIII. | Post digestion spike | NA | |
| XIV. | Compound quantitation LOQ/LOD/DL | Y | |
| XV. | Target compound identification | Y | |

VALIDATION CHECKLIST

SDG#: 280-197922

Date: 2/14/25

Laboratory: EETA

Reviewer: Kortney Curry

Method: 8015

| | Validation Area | Acceptable? Y/N/NA | Comments |
|-------|--|-----------------------|----------|
| I. | Case narrative | Y | |
| II. | Sample receipt/Technical holding times | Y | |
| III. | Instrument performance check/Tune | NA | |
| IV. | Initial calibration/ICV | Y | |
| V. | Continuing Calibration | Y | |
| VI. | Laboratory Blanks- MB | Y | |
| VI. | Field blanks | Y | |
| VII. | Surrogate spikes | Y | |
| VIII. | Matrix spike/Matrix spike duplicate | NA | |
| IX. | Laboratory control samples | Y | |
| X. | Field duplicates/Field triplicates | NA | |
| XI. | Internal standards | NA | |
| XII. | Compound quantitation LOQ/LOD/DL | Y | |
| XIII. | Target compound identification | Y | |

Attachment C

ADR Summary Report



Data Review Sample Summary Report by Analysis Method

Reviewed By: KAC (1/14/2025)

Approved By:

Laboratory: TAL DEN

| Client Sample ID | Lab Sample ID | Matrix | Sample Type | Preparation Method | Collection Date | Validation Code |
|------------------|---------------|--------|-------------|--------------------|-----------------|-----------------|
|------------------|---------------|--------|-------------|--------------------|-----------------|-----------------|

Lab Reporting Batch: 280-197922-1

Method: 365.1

| | | | | | | |
|----------------|-----------------|-------|------------------------|----------|-----------------------|-------|
| TMW38102024 | 280-197922-4 | Water | Field_Sample | Gen Prep | 10/10/2024 8:08:00 AM | S2AVE |
| TMW32102024 | 280-197922-3 | Water | Field_Sample | Gen Prep | 10/10/2024 9:55:00 AM | S2AVE |
| TMW21102024 | 280-197922-2 | Water | Field_Sample | Gen Prep | 10/10/2024 8:05:00 AM | S2AVE |
| TMW32102024MSD | 280-197922-3MSD | Water | Matrix_Spike_Duplicate | Gen Prep | 10/10/2024 9:55:00 AM | S2AVE |
| TMW32102024MS | 280-197922-3MS | Water | Matrix_Spike | Gen Prep | 10/10/2024 9:55:00 AM | S2AVE |

Method: 6020B

| | | | | | | |
|----------------|-----------------|-------|------------------------|-------|-----------------------|-------|
| TMW38102024 | 280-197922-7 | Water | Field_Sample | 3005A | 10/10/2024 8:08:00 AM | S2AVE |
| TMW21102024 | 280-197922-5 | Water | Field_Sample | 3005A | 10/10/2024 8:05:00 AM | S2AVE |
| TMW32102024 | 280-197922-6 | Water | Field_Sample | 3005A | 10/10/2024 9:55:00 AM | S2AVE |
| TMW32102024 | 280-197922-6 | Water | Field_Sample | 3020A | 10/10/2024 9:55:00 AM | S2AVE |
| TMW38102024 | 280-197922-7 | Water | Field_Sample | 3020A | 10/10/2024 8:08:00 AM | S2AVE |
| TMW38102024MSD | 280-197922-7MSD | Water | Matrix_Spike_Duplicate | 3020A | 10/10/2024 8:08:00 AM | S2AVE |
| TMW38102024MS | 280-197922-7MS | Water | Matrix_Spike | 3020A | 10/10/2024 8:08:00 AM | S2AVE |
| TMW21102024 | 280-197922-5 | Water | Field_Sample | 3020A | 10/10/2024 8:05:00 AM | S2AVE |

Method: 6850

| | | | | | | |
|-------------|--------------|-------|--------------|----------|-----------------------|-------|
| TMW21102024 | 280-197922-5 | Water | Field_Sample | Gen Prep | 10/10/2024 8:05:00 AM | S2AVE |
| TMW38102024 | 280-197922-7 | Water | Field_Sample | Gen Prep | 10/10/2024 8:08:00 AM | S2AVE |
| TMW32102024 | 280-197922-6 | Water | Field_Sample | Gen Prep | 10/10/2024 9:55:00 AM | S2AVE |

Method: 7470A

| | | | | | | |
|-------------|--------------|-------|--------------|-------|-----------------------|-------|
| TMW21102024 | 280-197922-5 | Water | Field_Sample | 7470A | 10/10/2024 8:05:00 AM | S2AVE |
| TMW38102024 | 280-197922-7 | Water | Field_Sample | 7470A | 10/10/2024 8:08:00 AM | S2AVE |
| TMW32102024 | 280-197922-6 | Water | Field_Sample | 7470A | 10/10/2024 9:55:00 AM | S2AVE |



Data Review Sample Summary Report by Analysis Method

Reviewed By: KAC (1/14/2025)

Approved By:

Laboratory: TAL DEN

| Client Sample ID | Lab Sample ID | Matrix | Sample Type | Preparation Method | Collection Date | Validation Code |
|--------------------------|-----------------|--------|------------------------|--------------------|-----------------------|-----------------|
| Method: 8015D-DRO | | | | | | |
| TMW21102024 | 280-197922-5 | Water | Field_Sample | 3510C | 10/10/2024 8:05:00 AM | S2AVE |
| Method: 8015D-GRO | | | | | | |
| QC10102024TB | 280-197922-1 | Water | Trip_Blank | Gen Prep | 10/10/2024 8:00:00 AM | S2AVE |
| TMW21102024 | 280-197922-5 | Water | Field_Sample | Gen Prep | 10/10/2024 8:05:00 AM | S2AVE |
| Method: 8260D | | | | | | |
| TMW32102024 | 280-197922-6 | Water | Field_Sample | 5030B | 10/10/2024 9:55:00 AM | S2AVE |
| TMW38102024 | 280-197922-7 | Water | Field_Sample | 5030B | 10/10/2024 8:08:00 AM | S2AVE |
| QC10102024TB | 280-197922-1 | Water | Trip_Blank | 5030B | 10/10/2024 8:00:00 AM | S2AVE |
| TMW21102024 | 280-197922-5 | Water | Field_Sample | 5030B | 10/10/2024 8:05:00 AM | S2AVE |
| Method: 8270E | | | | | | |
| TMW38102024 | 280-197922-7 | Water | Field_Sample | 3510C | 10/10/2024 8:08:00 AM | S2AVE |
| Method: 8330B | | | | | | |
| TMW32102024 | 280-197922-6 | Water | Field_Sample | 3535 | 10/10/2024 9:55:00 AM | S2AVE |
| TMW38102024 | 280-197922-7 | Water | Field_Sample | 3535 | 10/10/2024 8:08:00 AM | S2AVE |
| TMW21102024 | 280-197922-5 | Water | Field_Sample | 3535 | 10/10/2024 8:05:00 AM | S2AVE |
| Method: 9056A | | | | | | |
| TMW32102024DUP | 280-197922-3DUP | Water | Duplicate | Gen Prep | 10/10/2024 9:55:00 AM | S2AVE |
| TMW38102024 | 280-197922-4 | Water | Field_Sample | Gen Prep | 10/10/2024 8:08:00 AM | S2AVE |
| TMW32102024 | 280-197922-3 | Water | Field_Sample | Gen Prep | 10/10/2024 9:55:00 AM | S2AVE |
| TMW21102024 | 280-197922-2 | Water | Field_Sample | Gen Prep | 10/10/2024 8:05:00 AM | S2AVE |
| TMW32102024MS | 280-197922-3MS | Water | Matrix_Spike | Gen Prep | 10/10/2024 9:55:00 AM | S2AVE |
| TMW32102024MSD | 280-197922-3MSD | Water | Matrix_Spike_Duplicate | Gen Prep | 10/10/2024 9:55:00 AM | S2AVE |
| TMW21102024DUP | 280-197922-2DUP | Water | Duplicate | Gen Prep | 10/10/2024 8:05:00 AM | S2AVE |
| TMW21102024MS | 280-197922-2MS | Water | Matrix_Spike | Gen Prep | 10/10/2024 8:05:00 AM | S2AVE |



Data Review Sample Summary Report by Analysis Method

Reviewed By: KAC (1/14/2025)

Approved By:

Laboratory: TAL DEN

| <i>Client Sample ID</i> | <i>Lab Sample ID</i> | <i>Matrix</i> | <i>Sample Type</i> | <i>Preparation Method</i> | <i>Collection Date</i> | <i>Validation Code</i> |
|-------------------------|----------------------|---------------|------------------------|---------------------------|--------------------------|------------------------|
| Method: 9056A | | | | | | |
| TMW21102024MSD | 280-197922-2MSD | Water | Matrix_Spike_Duplicate | Gen Prep | 10/10/2024 8:05:00 AM | S2AVE |



Data Review Sample Summary Report by Analysis Method

Reviewed By: KAC (1/14/2025)

Approved By:

Laboratory: TAL DEN

| <i>Client Sample ID</i> | <i>Lab Sample ID</i> | <i>Matrix</i> | <i>Sample Type</i> | <i>Preparation Method</i> | <i>Collection Date</i> | <i>Validation Code</i> |
|-------------------------|----------------------|---------------|--------------------|---------------------------|------------------------|------------------------|
|-------------------------|----------------------|---------------|--------------------|---------------------------|------------------------|------------------------|

Validation Label Legend

| <i>Label Code</i> | <i>Label Description</i> | <i>EPA Level</i> |
|-------------------|---|--------------------------|
| S1VE | Stage_1_Validation_Electronic | N/A |
| S1VM | Stage_1_Validation_Manual | N/A |
| S1VEM | Stage_1_Validation_Electronic_and_Manual | N/A |
| S2AVE | Stage_2A_Validation_Electronic | Level 3 w/o calibration |
| S2AVM | Stage_2A_Validation_Manual | Level 3 w/o calibration |
| S2AVEM | Stage_2A_Validation_Electronic_and_Manual | Level 3 w/o calibration |
| S2BVE | Stage_2B_Validation_Electronic | Level 3 with calibration |
| S2BVM | Stage_2B_Validation_Manual | Level 3 with calibration |
| S2BVEM | Stage_2B_Validation_Electronic_and_Manual | Level 3 with calibration |
| S3VE | Stage_3_Validation_Electronic | Level 4 |
| S3VM | Stage_3_Validation_Manual | Level 4 |
| S3VEM | Stage_3_Validation_Electronic_and_Manual | Level 4 |
| S4VE | Stage_4_Validation_Electronic | Level 4 |
| S4VM | Stage_4_Validation_Manual | Level 4 |
| S4VEM | Stage_4_Validation_Electronic_and_Manual | Level 4 |
| NV | Not_Validated | N/A |



Data Review Summary

Lab Reporting Batch ID: 280-197922-1

Laboratory: TAL DEN

EDD Filename: 280-197922-1_52_2a_ParsonsFtWingate

eQAPP Name: Fort_Wingate_rev2_hexchrom_AQ

Validation Area

Note

| | |
|---|----|
| Technical Holding Times | SR |
| Temperature | A |
| Initial Calibration | N |
| Continuing Calibration/Initial Calibration Verification | N |
| Method Blanks | SR |
| Surrogate/Tracer Spikes | A |
| Matrix Spike/Matrix Spike Duplicates | SR |
| Laboratory Duplicates | SR |
| Laboratory Replicates | N |
| Laboratory Control Samples | A |
| Compound Quantitation | SR |
| Field Duplicates | N |
| Field Triplicates | N |
| Field Blanks | A |

A = Acceptable, N = Not provided/applicable, SR = See report

The contents of this report reflect findings made by ADR during Automated Data Review, manual applied qualifiers are not considered. Please refer to the Overall Qualifier Summary report for manual qualifiers.

Temperature Outliers

Lab Reporting Batch ID: 280-197922-1

Laboratory: TAL DEN

EDD Filename: 280-197922-1_52_2a_ParsonsFtWingate

eQAPP Name: Fort_Wingate_rev2_hexchrom_AQ

No Data Review Qualifiers Applied

QC Outlier Report: HoldingTimes

Lab Reporting Batch ID: 280-197922-1

Laboratory: TAL DEN

EDD Filename: 280-197922-1_52_2a_ParsonsFtWingate

eQAPP Name: Fort_Wingate_rev2_hexchrom_AQ HT_24hr

Method: 9056A

Preparation Method: Gen Prep

Matrix: Water

| Sample ID | Type | Actual | Criteria | Units | Flag |
|---------------------------|----------------------|--------|----------|-------|--|
| TMW32102024 (Initial/TOT) | Sampling To Analysis | 59.00 | 48.00 | HOURS | J- (all detects) UJ (all non-detects) |

Project Name and Number: Fort Wingate Depot

1/14/2025 1:52:36 PM

ADR version 1.9.0.325

Page 1 of 1

Trip Blank Outlier Report

Lab Reporting Batch ID: 280-197922-1

Laboratory: TAL DEN

EDD Filename: 280-197922-1_52_2a_ParsonsFtWingate

eQAPP Name: Fort_Wingate_rev2_hexchrom_AQ

No Data Review Qualifiers Applied

Method Blank Outlier Report

Lab Reporting Batch ID: 280-197922-1

Laboratory: TAL DEN

EDD Filename: 280-197922-1_52_2a_ParsonsFtWingate

eQAPP Name: Fort_Wingate_rev2_hexchrom_AQ HT_24hr

| | |
|---------|--------|
| Method: | 6020B |
| Method: | Method |

| Method Blank Sample ID | Analysis Date | Analyte | Result | Associated Samples |
|------------------------|------------------------|-------------------|------------------------|---|
| MB 280-670743/1-A | 10/22/2024 4:09:18 AM | ALUMINUM | 9.36 ug/L | TMW21102024 TMW32102024 TMW38102024 |
| MB 280-670767/1-A | 10/17/2024 12:44:23 PM | IRON MAGNESIUM | 28.5 ug/L 29.3 ug/L | TMW21102024 TMW32102024 TMW38102024 |

The following samples and their listed target analytes were qualified due to contamination reported in this blank

| Sample ID | Analyte | Reported Result | Modified Final Result |
|--------------------------|----------|-----------------|-----------------------|
| TMW21102024(Initial/DIS) | ALUMINUM | 12 ug/L | 200U ug/L |
| TMW21102024(Initial/TOT) | IRON | 37 ug/L | 200U ug/L |
| TMW32102024(Initial/DIS) | ALUMINUM | 13 ug/L | 200U ug/L |
| TMW32102024(Initial/TOT) | IRON | 29 ug/L | 200U ug/L |
| TMW38102024(Initial/DIS) | ALUMINUM | 16 ug/L | 200U ug/L |
| TMW38102024(Initial/TOT) | IRON | 120 ug/L | 200U ug/L |

Project Name and Number: Fort Wingate Depot Activity Northern Area - NM6213820974

1/14/2025 1:52:48 PM

ADR version 1.9.0.325

Page 1 of 1

Surrogate Outlier Report

Lab Reporting Batch ID: 280-197922-1

Laboratory: TAL DEN

EDD Filename: 280-197922-1_52_2a_ParsonsFtWingate

eQAPP Name: Fort_Wingate_rev2_hexchrom_AQ

No Data Review Qualifiers Applied

Lab Control Spike/Lab Control Spike Duplicate Outlier Report

Lab Reporting Batch ID: 280-197922-1

Laboratory: TAL DEN

EDD Filename: 280-197922-1_52_2a_ParsonsFtWingate

eQAPP Name: Fort_Wingate_rev2_hexchrom_AQ

No Data Review Qualifiers Applied

Matrix Spike/Matrix Spike Duplicate Outlier Report

Lab Reporting Batch ID: 280-197922-1

Laboratory: TAL DEN

EDD Filename: 280-197922-1_52_2a_ParsonsFtWingate

eQAPP Name: Fort_Wingate_rev2_hexchrom_AQ

Method: 6020B

| QC Sample ID (Associated Samples) | Compound | MS %R | MSD %R | %R Limits | RPD (Limits) | Affected Compounds | Flag |
|--|----------|----------|-----------|--------------|-----------------|-----------------------|---|
| TMW38102024MS (Total) TMW38102024MSD (Total) (TMW38102024) | SODIUM | 556 | -2656 | 85.00-117.00 | - | SODIUM | J (all detects) UJ (all non-detects) |
| TMW38102024MS (Total) TMW38102024MSD (Total) (TMW38102024) | CALCIUM | 73 | 16 | 87.00-118.00 | - | CALCIUM | J-(all detects) UJ(all non-detects) |

Sample concentrations are greater than 4 times the MS/MSD spike concentrations.
RECs could not be evaluated, and qualification was not warranted.

Lab Duplicate Outlier Report

Lab Reporting Batch ID: 280-197922-1

Laboratory: TAL DEN

EDD Filename:

eQAPP Name: Fort_Wingate_rev2_hexchrom_AQ

280-197922-1_52_2a_ParsonsFtWingate

Method: 9056A

Matrix: Water

| QC Sample ID (Associated Sample ID) | Analyte | Sample RPD | eQAPP RPD | Flag |
|---|----------|---------------|--------------|---|
| TMW32102024DUP (TMW32102024) | FLUORIDE | 29 | 10.00 | J (all detects) UJ (all non-detects) |

Project Name and Number: Fort Wingate Depot Activity Northern Area - NM6213820974

1/14/2025 1:54:04 PM

ADR version 1.9.0.325

Page 1 of 1

Reporting Limit Outliers

Lab Reporting Batch ID: 280-197922-1

Laboratory: TAL DEN

EDD Filename: 280-197922-1_52_2a_ParsonsFtWingate

eQAPP Name: Fort_Wingate_rev2_hexchrom_AQ HT_24hr

Method: 365.1

| SampleID | Analyte | Lab Qual | Result | Reporting Limit | RL Type | Units | Flag |
|-------------|---------------------|----------|--------|-----------------|---------|-------|-----------------|
| TMW21102024 | Orthophosphate as P | J | 38 | 50 | LOQ | ug/L | J (all detects) |
| TMW38102024 | Orthophosphate as P | J | 39 | 50 | LOQ | ug/L | J (all detects) |

Method: 6020B

| SampleID | Analyte | Lab Qual | Result | Reporting Limit | RL Type | Units | Flag |
|-------------|-----------|----------|--------|-----------------|---------|-------|-----------------|
| TMW21102024 | ALUMINUM | J | 46 | 200 | LOQ | ug/L | J (all detects) |
| | ARSENIC | J | 0.93 | 5.0 | LOQ | ug/L | |
| | CHROMIUM | J | 1.3 | 3.0 | LOQ | ug/L | |
| | COPPER | J | 1.5 | 2.0 | LOQ | ug/L | |
| | IRON | J | 37 | 200 | LOQ | ug/L | |
| | MANGANESE | J | 0.71 | 3.0 | LOQ | ug/L | |
| | NICKEL | J | 2.5 | 3.0 | LOQ | ug/L | |
| | POTASSIUM | J | 430 | 1000 | LOQ | ug/L | |
| | VANADIUM | J | 3.2 | 5.0 | LOQ | ug/L | |
| | ZINC | J | 4.4 | 10 | LOQ | ug/L | |
| TMW32102024 | ALUMINUM | J | 25 | 200 | LOQ | ug/L | J (all detects) |
| | ANTIMONY | J | 0.45 | 2.0 | LOQ | ug/L | |
| | ARSENIC | J | 3.5 | 5.0 | LOQ | ug/L | |
| | CADMIUM | J | 0.22 | 1.0 | LOQ | ug/L | |
| | CHROMIUM | J | 0.65 | 3.0 | LOQ | ug/L | |
| | COPPER | J | 0.86 | 2.0 | LOQ | ug/L | |
| | IRON | J | 29 | 200 | LOQ | ug/L | |
| | POTASSIUM | J | 760 | 1000 | LOQ | ug/L | |
| | SELENIUM | J | 2.8 | 5.0 | LOQ | ug/L | |
| | SILVER | J | 0.10 | 1.0 | LOQ | ug/L | |
| | VANADIUM | J | 2.1 | 5.0 | LOQ | ug/L | |
| | ZINC | J | 4.8 | 10 | LOQ | ug/L | |
| TMW38102024 | ALUMINUM | J | 18 | 200 | LOQ | ug/L | J (all detects) |
| | ARSENIC | J | 0.85 | 5.0 | LOQ | ug/L | |
| | CHROMIUM | J | 0.61 | 3.0 | LOQ | ug/L | |
| | COPPER | J | 1.1 | 2.0 | LOQ | ug/L | |
| | IRON | J | 120 | 200 | LOQ | ug/L | |
| | ZINC | J | 2.0 | 10 | LOQ | ug/L | |

Method: 9056A

| SampleID | Analyte | Lab Qual | Result | Reporting Limit | RL Type | Units | Flag |
|-------------|----------|----------|--------|-----------------|---------|-------|-----------------|
| TMW32102024 | FLUORIDE | J M | 940 | 1000 | LOQ | ug/L | J (all detects) |

Project Name and Number: Fort Wingate Depot Activity Northern Area - NM6213820974

1/14/2025 1:54:18 PM

ADR version 1.9.0.325

Page 1 of 1



Field QC Assignments and Associated Samples

EDD File Name: 280-197922-1

eQapp Name: Fort_Wingate_rev2_hexchrom_AQ HT_24hr

| | Associated Samples | Sample Collection Date |
|---|--------------------|------------------------|
| <div><div>Field QC QC10102024TB</div><div>QC Type: Trip_Blank</div></div> | | |
| | TMW21102024 | 10/10/2024 8:05:00 AM |
| | TMW38102024 | 10/10/2024 8:08:00 AM |
| | TMW32102024 | 10/10/2024 9:55:00 AM |



Data Qualifier Summary

Lab Reporting Batch ID: 280-197922-1

Laboratory: TAL DEN

EDD Filename:
280-197922-1_52_2a_ParsonsFtWingate_rev_rev

eQAPP Name: Fort_Wingate_rev2_hexchrom_AQ HT_24hr

Method Category: GENCHEM

| Sample ID:TMW21102024 | | 10/10/2024 8:05:00 Collected:AM | | Analysis Type:Initial/TOT | | | | Dilution: 1 | |
|-----------------------|------------|------------------------------------|----|---------------------------|----|---------|-------|------------------|-------------|
| Analyte | Lab Result | Lab Qual | DL | DL Type | RL | RL Type | Units | Data Review Qual | Reason Code |
| Orthophosphate as P | 38 | J | 40 | LOD | 50 | LOQ | ug/L | J | TR |

| Sample ID:TMW38102024 | | 10/10/2024 8:08:00 Collected:AM | | Analysis Type:Initial/TOT | | | | Dilution: 1 | |
|-----------------------|------------|------------------------------------|----|---------------------------|----|---------|-------|------------------|-------------|
| Analyte | Lab Result | Lab Qual | DL | DL Type | RL | RL Type | Units | Data Review Qual | Reason Code |
| Orthophosphate as P | 39 | J | 40 | LOD | 50 | LOQ | ug/L | J | TR |

Method Category: GENCHEM

| Sample ID:TMW32102024 | | 10/10/2024 9:55:00 Collected:AM | | Analysis Type:Initial/TOT | | | | Dilution: 1 | |
|-----------------------|------------|------------------------------------|-----|---------------------------|------|---------|-------|------------------|-------------|
| Analyte | Lab Result | Lab Qual | DL | DL Type | RL | RL Type | Units | Data Review Qual | Reason Code |
| FLUORIDE | 940 | J M | 500 | LOD | 1000 | LOQ | ug/L | J | TR, DU2 |
| Nitrate as N | 3000 | H | 200 | LOD | 500 | LOQ | ug/L | J- | SC1 |

Method Category: METALS

| Sample ID:TMW21102024 | | 10/10/2024 8:05:00 Collected:AM | | Analysis Type:Initial/DIS | | | | Dilution: 1 | |
|-----------------------|------------|------------------------------------|-----|---------------------------|------|---------|-------|------------------|-------------|
| Analyte | Lab Result | Lab Qual | DL | DL Type | RL | RL Type | Units | Data Review Qual | Reason Code |
| ALUMINUM | 12 | J | 30 | LOD | 200 | LOQ | ug/L | U | BLT/BLU |
| ARSENIC | 0.95 | J | 2.0 | LOD | 5.0 | LOQ | ug/L | J | TR |
| CHROMIUM | 1.3 | J | 1.8 | LOD | 3.0 | LOQ | ug/L | J | TR |
| COPPER | 1.4 | J | 1.8 | LOD | 2.0 | LOQ | ug/L | J | TR |
| IRON | 19 | J | 40 | LOD | 200 | LOQ | ug/L | J | TR |
| POTASSIUM | 450 | J | 76 | LOD | 1000 | LOQ | ug/L | J | TR |
| VANADIUM | 2.9 | J | 3.0 | LOD | 5.0 | LOQ | ug/L | J | TR |
| ZINC | 4.4 | J | 8.0 | LOD | 10 | LOQ | ug/L | J | TR |

* denotes a non-reportable result

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2/11/2025 3:07:33 PM

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Page 1 of 7



Data Qualifier Summary

Lab Reporting Batch ID: 280-197922-1

Laboratory: TAL DEN

EDD Filename:
280-197922-1_52_2a_ParsonsFtWingate_rev_rev

eQAPP Name: Fort_Wingate_rev2_hexchrom_AQ HT_24hr

Method Category: METALS

| Sample ID:TMW21102024 | | 10/10/2024 8:05:00 Collected: AM | | Analysis Type: Initial/TOT | | | | Dilution: 1 | |
|-----------------------|------------|-------------------------------------|-----|----------------------------|------|---------|-------|------------------|-------------|
| Analyte | Lab Result | Lab Qual | DL | DL Type | RL | RL Type | Units | Data Review Qual | Reason Code |
| ALUMINUM | 46 | J | 30 | LOD | 200 | LOQ | ug/L | J | TR |
| ARSENIC | 0.93 | J | 2.0 | LOD | 5.0 | LOQ | ug/L | J | TR |
| CHROMIUM | 1.3 | J | 1.8 | LOD | 3.0 | LOQ | ug/L | J | TR |
| COPPER | 1.5 | J | 1.8 | LOD | 2.0 | LOQ | ug/L | J | TR |
| IRON | 37 | J | 40 | LOD | 200 | LOQ | ug/L | U | BLT/BLU |
| MANGANESE | 0.71 | J | 1.8 | LOD | 3.0 | LOQ | ug/L | J | TR |
| NICKEL | 2.5 | J | 1.9 | LOD | 3.0 | LOQ | ug/L | J | TR |
| POTASSIUM | 430 | J | 76 | LOD | 1000 | LOQ | ug/L | J | TR |
| VANADIUM | 3.2 | J | 3.0 | LOD | 5.0 | LOQ | ug/L | J | TR |

| Sample ID:TMW32102024 | | 10/10/2024 9:55:00 Collected: AM | | Analysis Type: Initial/DIS | | | | Dilution: 1 | |
|-----------------------|------------|-------------------------------------|-----|----------------------------|------|---------|-------|------------------|-------------|
| Analyte | Lab Result | Lab Qual | DL | DL Type | RL | RL Type | Units | Data Review Qual | Reason Code |
| ALUMINUM | 13 | J | 30 | LOD | 200 | LOQ | ug/L | U | BLT/BLU |
| ANTIMONY | 0.40 | J | 1.0 | LOD | 2.0 | LOQ | ug/L | J | TR |
| ARSENIC | 1.4 | J | 2.0 | LOD | 5.0 | LOQ | ug/L | J | TR |
| IRON | 8.8 | J | 40 | LOD | 200 | LOQ | ug/L | J | TR |
| POTASSIUM | 860 | J | 76 | LOD | 1000 | LOQ | ug/L | J | TR |
| SELENIUM | 3.1 | J | 4.0 | LOD | 5.0 | LOQ | ug/L | J | TR |
| VANADIUM | 2.2 | J | 3.0 | LOD | 5.0 | LOQ | ug/L | J | TR |

| Sample ID:TMW32102024 | | 10/10/2024 9:55:00 Collected: AM | | Analysis Type: Initial/TOT | | | | Dilution: 1 | |
|-----------------------|------------|-------------------------------------|------|----------------------------|------|---------|-------|------------------|-------------|
| Analyte | Lab Result | Lab Qual | DL | DL Type | RL | RL Type | Units | Data Review Qual | Reason Code |
| ALUMINUM | 25 | J | 30 | LOD | 200 | LOQ | ug/L | J | TR |
| ANTIMONY | 0.45 | J | 1.0 | LOD | 2.0 | LOQ | ug/L | J | TR |
| ARSENIC | 3.5 | J | 2.0 | LOD | 5.0 | LOQ | ug/L | J | TR |
| CADMIUM | 0.22 | J | 0.75 | LOD | 1.0 | LOQ | ug/L | J | TR |
| CHROMIUM | 0.65 | J | 1.8 | LOD | 3.0 | LOQ | ug/L | J | TR |
| COPPER | 0.86 | J | 1.8 | LOD | 2.0 | LOQ | ug/L | J | TR |
| IRON | 29 | J | 40 | LOD | 200 | LOQ | ug/L | U | BLT/BLU |
| POTASSIUM | 760 | J | 76 | LOD | 1000 | LOQ | ug/L | J | TR |

* denotes a non-reportable result

Project Name and Number: Fort Wingate Depot Activity Northern Area - USACE Project: USACE Project: NM6213820974

2/11/2025 3:07:33 PM

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Page 2 of 7



Data Qualifier Summary

Lab Reporting Batch ID: 280-197922-1

Laboratory: TAL DEN

EDD Filename:
280-197922-1_52_2a_ParsonsFtWingate_rev_rev

eQAPP Name: Fort_Wingate_rev2_hexchrom_AQ HT_24hr

Method Category: METALS

| Sample ID:TMW32102024 | | 10/10/2024 9:55:00 Collected: AM | | Analysis Type: Initial/TOT | | | | Dilution: 1 | |
|-----------------------|------------|-------------------------------------|------|----------------------------|-----|---------|-------|------------------|-------------|
| Analyte | Lab Result | Lab Qual | DL | DL Type | RL | RL Type | Units | Data Review Qual | Reason Code |
| SELENIUM | 2.8 | J | 4.0 | LOD | 5.0 | LOQ | ug/L | J | TR |
| SILVER | 0.10 | J | 0.15 | LOD | 1.0 | LOQ | ug/L | U | ICB/CCB |
| VANADIUM | 2.1 | J | 3.0 | LOD | 5.0 | LOQ | ug/L | J | TR |
| ZINC | 4.8 | J | 8.0 | LOD | 10 | LOQ | ug/L | J | TR |

| Sample ID:TMW38102024 | | 10/10/2024 8:08:00 Collected: AM | | Analysis Type: Initial/DIS | | | | Dilution: 1 | |
|-----------------------|------------|-------------------------------------|-----|----------------------------|-----|---------|-------|------------------|-------------|
| Analyte | Lab Result | Lab Qual | DL | DL Type | RL | RL Type | Units | Data Review Qual | Reason Code |
| ALUMINUM | 16 | J | 30 | LOD | 200 | LOQ | ug/L | U | BLT/BLU |
| ARSENIC | 0.66 | J | 2.0 | LOD | 5.0 | LOQ | ug/L | J | TR |
| IRON | 79 | J | 40 | LOD | 200 | LOQ | ug/L | J | TR |

| Sample ID:TMW38102024 | | 10/10/2024 8:08:00 Collected: AM | | Analysis Type: Initial/TOT | | | | Dilution: 1 | |
|-----------------------|------------|-------------------------------------|-----|----------------------------|-----|---------|-------|------------------|-------------|
| Analyte | Lab Result | Lab Qual | DL | DL Type | RL | RL Type | Units | Data Review Qual | Reason Code |
| ALUMINUM | 18 | J | 30 | LOD | 200 | LOQ | ug/L | J | TR |
| ARSENIC | 0.85 | J | 2.0 | LOD | 5.0 | LOQ | ug/L | J | TR |
| CHROMIUM | 0.61 | J | 1.8 | LOD | 3.0 | LOQ | ug/L | J | TR |
| COPPER | 1.1 | J | 1.8 | LOD | 2.0 | LOQ | ug/L | J | TR |
| IRON | 120 | J | 40 | LOD | 200 | LOQ | ug/L | U | BLT/BLU |
| ZINC | 2.0 | J | 8.0 | LOD | 10 | LOQ | ug/L | J | TR |

Method Category: VOA

| Sample ID:QC10102024TB | | 10/10/2024 8:00:00 Collected: AM | | Analysis Type: Initial/TOT | | | | Dilution: 1 | |
|---------------------------|------------|-------------------------------------|------|----------------------------|-----|---------|-------|------------------|-------------|
| Analyte | Lab Result | Lab Qual | DL | DL Type | RL | RL Type | Units | Data Review Qual | Reason Code |
| 1,1,1,2-TETRACHLOROETHANE | 0.80 | U | 0.80 | LOD | 1.0 | LOQ | ug/L | UJ | PR1 |
| 1,1,1-TRICHLOROETHANE | 0.50 | U | 0.50 | LOD | 1.0 | LOQ | ug/L | UJ | PR1 |
| 1,1,2,2-TETRACHLOROETHANE | 0.80 | U | 0.80 | LOD | 1.0 | LOQ | ug/L | UJ | PR1 |
| 1,1,2-TRICHLOROETHANE | 0.80 | U | 0.80 | LOD | 1.0 | LOQ | ug/L | UJ | PR1 |

* denotes a non-reportable result

Project Name and Number: Fort Wingate Depot Activity Northern Area - USACE Project: USACE Project: NM6213820974

2/11/2025 3:07:33 PM

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Page 3 of 7



Data Qualifier Summary

Lab Reporting Batch ID: 280-197922-1

Laboratory: TAL DEN

EDD Filename:
280-197922-1_52_2a_ParsonsFtWingate_rev_rev

eQAPP Name: Fort_Wingate_rev2_hexchrom_AQ HT_24hr

Method Category: VOA

| 10/10/2024 8:00:00 | | | | | | | | | |
|-----------------------------|------------|---------------|------|----------------------------|-----|---------|-------|------------------|-------------|
| Sample ID: QC10102024TB | | Collected: AM | | Analysis Type: Initial/TOT | | | | Dilution: 1 | |
| Analyte | Lab Result | Lab Qual | DL | DL Type | RL | RL Type | Units | Data Review Qual | Reason Code |
| 1,1-DICHLOROETHANE | 0.80 | U | 0.80 | LOD | 1.0 | LOQ | ug/L | UJ | PR1 |
| 1,1-DICHLOROETHENE | 0.80 | U | 0.80 | LOD | 1.0 | LOQ | ug/L | UJ | PR1 |
| 1,1-DICHLOROPROPENE | 0.80 | U | 0.80 | LOD | 1.0 | LOQ | ug/L | UJ | PR1 |
| 1,2,3-TRICHLOROBENZENE | 2.5 | U | 2.5 | LOD | 4.0 | LOQ | ug/L | UJ | PR1 |
| 1,2,3-TRICHLOROPROPANE | 1.8 | U | 1.8 | LOD | 2.5 | LOQ | ug/L | UJ | PR1 |
| 1,2,4-TRICHLOROBENZENE | 0.80 | U | 0.80 | LOD | 1.0 | LOQ | ug/L | UJ | PR1 |
| 1,2,4-TRIMETHYLBENZENE | 0.40 | U | 0.40 | LOD | 1.0 | LOQ | ug/L | UJ | PR1 |
| 1,2-DIBROMO-3-CHLOROPROPANE | 4.0 | U | 4.0 | LOD | 5.0 | LOQ | ug/L | UJ | PR1 |
| 1,2-DIBROMOETHANE | 0.80 | U | 0.80 | LOD | 1.0 | LOQ | ug/L | UJ | PR1 |
| 1,2-DICHLOROBENZENE | 0.50 | U | 0.50 | LOD | 1.0 | LOQ | ug/L | UJ | PR1 |
| 1,2-DICHLOROETHANE | 0.50 | U | 0.50 | LOD | 1.0 | LOQ | ug/L | UJ | PR1 |
| 1,2-DICHLOROPROPANE | 0.80 | U | 0.80 | LOD | 1.0 | LOQ | ug/L | UJ | PR1 |
| 1,3,5-TRIMETHYLBENZENE | 0.50 | U | 0.50 | LOD | 1.0 | LOQ | ug/L | UJ | PR1 |
| 1,3-DICHLOROBENZENE | 0.40 | U | 0.40 | LOD | 1.0 | LOQ | ug/L | UJ | PR1 |
| 1,3-DICHLOROPROPANE | 0.80 | U | 0.80 | LOD | 1.0 | LOQ | ug/L | UJ | PR1 |
| 1,4-DICHLOROBENZENE | 0.50 | U | 0.50 | LOD | 1.0 | LOQ | ug/L | UJ | PR1 |
| 2,2-DICHLOROPROPANE | 0.80 | U | 0.80 | LOD | 1.0 | LOQ | ug/L | UJ | PR1 |
| 2-Butanone [MEK] | 8.0 | U | 8.0 | LOD | 10 | LOQ | ug/L | UJ | PR1 |
| 2-CHLOROTOLUENE | 0.40 | U | 0.40 | LOD | 1.0 | LOQ | ug/L | UJ | PR1 |
| 2-HEXANONE | 4.0 | U | 4.0 | LOD | 5.0 | LOQ | ug/L | UJ | PR1 |
| 4-CHLOROTOLUENE | 0.80 | U | 0.80 | LOD | 1.0 | LOQ | ug/L | UJ | PR1 |
| 4-ISOPROPYLTOLUENE | 0.80 | U | 0.80 | LOD | 1.0 | LOQ | ug/L | UJ | PR1 |
| 4-Methyl-2-pentanone [MIBK] | 3.2 | U | 3.2 | LOD | 5.0 | LOQ | ug/L | UJ | PR1 |
| ACETONE | 8.0 | U | 8.0 | LOD | 15 | LOQ | ug/L | UJ | PR1 |
| BENZENE | 0.80 | U | 0.80 | LOD | 1.0 | LOQ | ug/L | UJ | PR1 |
| BROMOBENZENE | 0.50 | U | 0.50 | LOD | 1.0 | LOQ | ug/L | UJ | PR1 |
| BROMOCHLOROMETHANE | 0.80 | U | 0.80 | LOD | 1.0 | LOQ | ug/L | UJ | PR1 |
| BROMODICHLOROMETHANE | 0.50 | U | 0.50 | LOD | 1.0 | LOQ | ug/L | UJ | PR1 |
| BROMOFORM | 1.8 | U | 1.8 | LOD | 2.0 | LOQ | ug/L | UJ | PR1 |
| BROMOMETHANE | 4.0 | U | 4.0 | LOD | 5.0 | LOQ | ug/L | UJ | PR1 |
| CARBON DISULFIDE | 0.80 | U | 0.80 | LOD | 2.0 | LOQ | ug/L | UJ | PR1 |

* denotes a non-reportable result

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2/11/2025 3:07:33 PM

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Page 4 of 7



Data Qualifier Summary

Lab Reporting Batch ID: 280-197922-1

Laboratory: TAL DEN

EDD Filename:
280-197922-1_52_2a_ParsonsFtWingate_rev_rev

eQAPP Name: Fort_Wingate_rev2_hexchrom_AQ HT_24hr

Method Category: VOA

| 10/10/2024 8:00:00 | | | | | | | | | |
|---------------------------|------------|--------------|------|---------------------------|-----|---------|-------|------------------|-------------|
| Sample ID:QC10102024TB | | Collected:AM | | Analysis Type:Initial/TOT | | | | Dilution: 1 | |
| Analyte | Lab Result | Lab Qual | DL | DL Type | RL | RL Type | Units | Data Review Qual | Reason Code |
| CARBON TETRACHLORIDE | 0.50 | U | 0.50 | LOD | 1.0 | LOQ | ug/L | UJ | PR1 |
| CHLOROBENZENE | 0.80 | U | 0.80 | LOD | 1.0 | LOQ | ug/L | UJ | PR1 |
| CHLOROETHANE | 1.0 | U | 1.0 | LOD | 2.0 | LOQ | ug/L | UJ | PR1 |
| CHLOROFORM | 0.80 | U | 0.80 | LOD | 1.0 | LOQ | ug/L | UJ | PR1 |
| CHLOROMETHANE | 1.0 | U | 1.0 | LOD | 2.0 | LOQ | ug/L | UJ | PR1 |
| CIS-1,2-DICHLOROETHENE | 0.40 | U | 0.40 | LOD | 1.0 | LOQ | ug/L | UJ | PR1 |
| CIS-1,3-DICHLOROPROPENE | 0.40 | U | 0.40 | LOD | 1.0 | LOQ | ug/L | UJ | PR1 |
| DIBROMOCHLOROMETHANE | 0.50 | U | 0.50 | LOD | 1.0 | LOQ | ug/L | UJ | PR1 |
| DIBROMOMETHANE | 0.40 | U | 0.40 | LOD | 1.0 | LOQ | ug/L | UJ | PR1 |
| DICHLORODIFLUOROMETHANE | 1.0 | U | 1.0 | LOD | 2.0 | LOQ | ug/L | UJ | PR1 |
| ETHYLBENZENE | 0.40 | U | 0.40 | LOD | 1.0 | LOQ | ug/L | UJ | PR1 |
| HEXACHLOROBUTADIENE | 1.8 | U | 1.8 | LOD | 2.0 | LOQ | ug/L | UJ | PR1 |
| ISOPROPYLBENZENE | 0.50 | U | 0.50 | LOD | 1.0 | LOQ | ug/L | UJ | PR1 |
| METHYL ACETATE | 4.0 | U | 4.0 | LOD | 5.0 | LOQ | ug/L | UJ | PR1 |
| METHYL TERT-BUTYL ETHER | 0.80 | U | 0.80 | LOD | 5.0 | LOQ | ug/L | UJ | PR1 |
| METHYLENE CHLORIDE | 1.8 | U | 1.8 | LOD | 2.0 | LOQ | ug/L | UJ | PR1 |
| m-Xylene & p-Xylene | 0.80 | U | 0.80 | LOD | 2.0 | LOQ | ug/L | UJ | PR1 |
| NAPHTHALENE | 2.0 | U | 2.0 | LOD | 3.0 | LOQ | ug/L | UJ | PR1 |
| N-BUTYLBENZENE | 0.80 | U | 0.80 | LOD | 1.0 | LOQ | ug/L | UJ | PR1 |
| N-PROPYLBENZENE | 0.80 | U | 0.80 | LOD | 1.0 | LOQ | ug/L | UJ | PR1 |
| O-XYLENE | 0.40 | U | 0.40 | LOD | 1.0 | LOQ | ug/L | UJ | PR1 |
| SEC-BUTYLBENZENE | 0.80 | U | 0.80 | LOD | 1.0 | LOQ | ug/L | UJ | PR1 |
| STYRENE | 0.80 | U | 0.80 | LOD | 1.0 | LOQ | ug/L | UJ | PR1 |
| TERT-BUTYLBENZENE | 0.80 | U | 0.80 | LOD | 1.0 | LOQ | ug/L | UJ | PR1 |
| TETRACHLOROETHENE | 0.80 | U | 0.80 | LOD | 1.0 | LOQ | ug/L | UJ | PR1 |
| TOLUENE | 0.40 | U | 0.40 | LOD | 1.0 | LOQ | ug/L | UJ | PR1 |
| TRANS-1,2-DICHLOROETHENE | 0.50 | U | 0.50 | LOD | 1.0 | LOQ | ug/L | UJ | PR1 |
| TRANS-1,3-DICHLOROPROPENE | 0.40 | U | 0.40 | LOD | 1.0 | LOQ | ug/L | UJ | PR1 |
| TRICHLOROETHENE | 0.40 | U | 0.40 | LOD | 1.0 | LOQ | ug/L | UJ | PR1 |
| TRICHLOROFLUOROMETHANE | 0.80 | U | 0.80 | LOD | 2.0 | LOQ | ug/L | UJ | PR1 |
| VINYL CHLORIDE | 0.40 | U | 0.40 | LOD | 1.0 | LOQ | ug/L | UJ | PR1 |

* denotes a non-reportable result

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Page 5 of 7



Data Qualifier Summary

Lab Reporting Batch ID: 280-197922-1

Laboratory: TAL DEN

EDD Filename:
280-197922-1_52_2a_ParsonsFtWingate_rev_rev

eQAPP Name: Fort_Wingate_rev2_hexchrom_AQ HT_24hr

Reason Code Legend

| <i>Reason Code</i> | <i>Description</i> |
|--------------------|---------------------------------|
| BLT/BLU | Method Blank Contamination |
| DU2 | Laboratory Duplicate Precision |
| ICB/CCB | Calibration Blank Contamination |
| MD1 | Matrix Spike Upper Estimation |
| MD2 | Matrix Spike Lower Estimation |
| PR1 | Preservation |
| SC1 | Sampling to Analysis Estimation |
| TR | Reporting Limit Trace Value |

* denotes a non-reportable result

Project Name and Number: Fort Wingate Depot Activity Northern Area - USACE Project: USACE Project: NM6213820974

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Page 7 of 7