

## Laboratory Data Consultants, Inc.

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Tronox, LLC  
P.O. Box 55  
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ATTN: Ms. Susan Crowley

December 29, 2010

SUBJECT: 2010 Annual Remedial Performance Sampling, Data Validation

Dear Ms. Crowley,

Enclosed are the final validation reports for the fractions listed below. These SDGs were received on October 20, 2010. Attachment 1 is a summary of the samples that were reviewed for each analysis.

**LDC Project # 24203:**

**SDG #**

**Fraction**

337662, 338619, 340066, 340161 340226, 340229, 340275, 340276 340278, 340887, 341684, 343130 343913, 339791, 339977	Chromium, Wet Chemistry
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The data validation was performed under Stage 2A guidelines. The analyses were validated using the following documents, as applicable to each method:

- USEPA, Contract Laboratory Program National Functional Guidelines for Inorganic Data Review, October 2004
- Region 9 Superfund Data Evaluation/Validation Guidance, NDEP Guidance, May 2006
- EPA SW 846, Third Edition, Test Methods for Evaluating Solid Waste, update 1, July 1992; update IIA, August 1993; update II, September 1994; update IIB, January 1995; update III, December 1996; update IIIA, April 1998; IIIB, November 2004; Update IV, February 2007

Please feel free to contact us if you have any questions.

Sincerely,

Erlinda T. Rauto  
Operations Manager/Senior Chemist



## Laboratory Data Consultants, Inc. Data Validation Report

**Project/Site Name:** 2010 Annual Remedial Performance Sampling  
**Collection Date:** August 3 through August 4, 2010  
**LDC Report Date:** December 19, 2010  
**Matrix:** Water  
**Parameters:** Chromium  
**Validation Level:** Stage 2A  
**Laboratory:** MWH Laboratories, Inc.  
**Sample Delivery Group (SDG):** 340066

### Sample Identification

M-99	M-88
M-25	I-V
M-92	VD080410
M-97	I-ADMS
M-14A	I-ADMSD
M-115	I-ACMS
M-17A	I-ACMSD
M-34	
M-35	
M-19	
M-39	
I-K	
I-J	
I-AD	
I-Z	
I-I	
I-AC	
M-68	
M-74	
M-73	

## Introduction

This data review covers 27 water samples listed on the cover sheet including dilutions and reanalysis as applicable. The analyses were per EPA SW 846 Method 6010 for Chromium.

This review follows a modified outline of the USEPA Contract Laboratory Program National Functional Guidelines for Inorganic Data Review (October 2004) and the EPA Region 9 Superfund Data Evaluation/Validation Guidance, NDEP guidance (May 2006).

A qualification summary table is provided at the end of this report if data has been qualified. Flags are classified as P (protocol) or A (advisory) to indicate whether the flag is due to a laboratory deviation from a specified protocol or is of technical advisory nature.

Raw data were not reviewed for this SDG. The review was based on QC data.

The following are definitions of the data qualifiers:

- J+ Data are qualified as estimated, with a high bias likely to occur. False positives or false negatives are unlikely to have been reported.
- J- Data are qualified as estimated, with a low bias likely to occur. False positives or false negatives are unlikely to have been reported.
- J Data are qualified as estimated; it is not possible to assess the direction of the potential bias. False positives or false negatives are unlikely to have been reported.
- U Indicates the compound or analyte was analyzed for but not detected at or above the stated limit.
- R Data are qualified as rejected. There is a significant potential for the reporting of false negatives or false positives.
- UJ Indicates the compound or analyte was analyzed for but not detected. The sample detection limit is an estimated value.
- A Indicates the finding is based upon technical validation criteria.
- P Indicates the finding is related to a protocol/contractual deviation.
- None Indicates the data was not significantly impacted by the finding, therefore qualification was not required.

## **I. Technical Holding Times**

All technical holding time requirements were met.

The chain-of-custodies were reviewed for documentation of cooler temperatures. All cooler temperatures met validation criteria.

## **II. ICPMS Tune**

ICP-MS was not utilized in this SDG.

## **III. Calibration**

Calibration data were not reviewed for Stage 2A.

## **IV. Blanks**

Method blanks were reviewed for each matrix as applicable. No chromium was found in the preparation blanks.

No field blanks were identified in this SDG.

## **V. ICP Interference Check Sample (ICS) Analysis**

ICP Interference check sample analysis data were not reviewed for Stage 2A.

## **VI. Matrix Spike Analysis**

Matrix spike (MS) and matrix spike duplicate (MSD) samples were reviewed for each matrix as applicable. Percent recoveries (%R) and relative percent differences (RPD) were within QC limits.

## **VII. Duplicate Sample Analysis**

Duplicate (DUP) sample analyses were reviewed for each matrix as applicable.

## **VIII. Laboratory Control Samples (LCS)**

Laboratory control samples were reviewed for each matrix as applicable. Percent recoveries (%R) and relative percent differences (RPD) were within QC limits.

## **IX. Internal Standards (ICP-MS)**

ICP-MS was not utilized in this SDG.

## **X. Furnace Atomic Absorption QC**

Graphite furnace atomic absorption was not utilized in this SDG.

## XI. ICP Serial Dilution

ICP serial dilution analysis data were not reviewed for Stage 2A.

## XII. Sample Result Verification

Raw data were not reviewed for this SDG.

## XIII. Overall Assessment of Data

Data flags are summarized at the end of this report if data has been qualified.

## XIV. Field Duplicates

Samples M-19 and VD080410 were identified as field duplicates. No chromium was detected in any of the samples with the following exceptions:

Analyte	Concentration (mg/L)		RPD (Limits)	Difference (Limits)	Flags	A or P
	M-19	VD080410				
Chromium	0.32	0.34	6 (≤30)	-	-	-

**2010 Annual Remedial Performance Sampling  
Chromium - Data Qualification Summary - SDG 340066**

No Sample Data Qualified in this SDG

**2010 Annual Remedial Performance Sampling  
Chromium - Laboratory Blank Data Qualification Summary - SDG 340066**

No Sample Data Qualified in this SDG

**2010 Annual Remedial Performance Sampling  
Chromium - Field Blank Data Qualification Summary - SDG 340066**

No Sample Data Qualified in this SDG

Tronox Northgate Henderson

LDC #: 24203C4  
 SDG #: 340066  
 Laboratory: MWH Laboratories

VALIDATION COMPLETENESS WORKSHEET

Stage 2  $\beta$  A

9/1/10

Date: 8-1-10  
 Page: 1 of 1  
 Reviewer: MG  
 2nd Reviewer: [Signature]

METHOD: Chromium (EPA SW 846 Method 6010)

The samples listed below were reviewed for each of the following validation areas. Validation findings are noted in attached validation findings worksheets.

	Validation Area		Comments
I.	Technical holding times	A	Sampling dates: 8-3-10 through 8-4-10
II.	ICP/MS Tune	N	not utilized
III.	Calibration	N	
IV.	Blanks	A	
V.	ICP Interference Check Sample (ICS) Analysis	N	
VI.	Matrix Spike Analysis	A	MS/MSD
VII.	Duplicate Sample Analysis	N	
VIII.	Laboratory Control Samples (LCS)	A	LCS/LCSD
IX.	Internal Standard (ICP-MS)	N	not utilized
X.	Furnace Atomic Absorption QC	N	" "
XI.	ICP Serial Dilution	N	not performed
XII.	Sample Result Verification	N	
XIII.	Overall Assessment of Data	A	
XIV.	Field Duplicates	SW	D = 10 + 23
XV.	Field Blanks	N	

Note: A = Acceptable  
 N = Not provided/applicable  
 SW = See worksheet

ND = No compounds detected  
 R = Rinsate  
 FB = Field blank

D = Duplicate  
 TB = Trip blank  
 EB = Equipment blank

Validated Samples:

all water

1	M-99	11	M-39	21	M-88	31	
2	M-25	12	I-K	22	I-V	32	
3	M-92	13	I-J	23	VD080410	33	
4	M-97	14	I-AD	24	I-ADMS	34	
5	M-14A	15	I-Z	25	I-ADMSD	35	
6	M-115	16	I-H	26	I-ACMS	36	
7	M-17A	17	I-AC	27	I-ACMSD	37	
8	M-34	18	M-68	28		38	
9	M-35	19	M-74	29	PBW1	39	
10	M-19	20	M-73	30	PBW2	40	

Notes: \_\_\_\_\_  
 \_\_\_\_\_  
 \_\_\_\_\_



LDC#: 24203C4

**VALIDATION FINDINGS WORKSHEET**  
**Field Duplicates**

Page: 1 of 1  
Reviewer: MG  
2nd Reviewer: [Signature]

**METHOD:** Metals (EPA Method 6010B/6020/7000)

- Y N NA Were field duplicate pairs identified in this SDG?
- Y N NA Were target analytes detected in the field duplicate pairs?

Analyte	Concentration (mg/L)		( $\leq 30$ )	(mg/L)	(mg/L)	Qualifications (Parent Only)
	10	23	RPD	Difference	Limits	
Chromium	0.32	0.34	6			

V:\FIELD DUPLICATES\FD\_inorganic\24203C4.wpd

## Laboratory Data Consultants, Inc. Data Validation Report

**Project/Site Name:** 2010 Annual Remedial Performance Sampling  
**Collection Date:** August 4, 2010  
**LDC Report Date:** December 22, 2010  
**Matrix:** Water  
**Parameters:** Chromium  
**Validation Level:** Stage 2A & 4  
**Laboratory:** MWH Laboratories, Inc.  
**Sample Delivery Group (SDG):** 340161

### Sample Identification

ART-1  
ART-2  
ART-3  
ART-4  
ART-6  
ART-7\*\*  
ART-8\*\*  
PC-99R2/R3\*\*  
PC-115R\*\*  
PC-116R\*\*  
PC-117\*\*  
PC-118\*\*  
PC-119  
PC-120  
PC-121  
PC-133  
ART-9  
PC-117MS  
PC-117MSD  
PC-118MS

\*\*Indicates sample underwent Stage 4 review

## Introduction

This data review covers 21 water samples listed on the cover sheet including dilutions and reanalysis as applicable. The analyses were per EPA SW 846 Method 6010 for Chromium.

This review follows a modified outline of the USEPA Contract Laboratory Program National Functional Guidelines for Inorganic Data Review (October 2004) and the EPA Region 9 Superfund Data Evaluation/Validation Guidance, NDEP guidance (May 2006).

A qualification summary table is provided at the end of this report if data has been qualified. Flags are classified as P (protocol) or A (advisory) to indicate whether the flag is due to a laboratory deviation from a specified protocol or is of technical advisory nature.

Samples indicated by a double asterisk on the front cover underwent a Stage 4 review. A Stage 2A review was performed on all of the other samples. Raw data were not evaluated for the samples reviewed by Stage 2A criteria since this review is based on QC data.

The following are definitions of the data qualifiers:

- J+ Data are qualified as estimated, with a high bias likely to occur. False positives or false negatives are unlikely to have been reported.
- J- Data are qualified as estimated, with a low bias likely to occur. False positives or false negatives are unlikely to have been reported.
- J Data are qualified as estimated; it is not possible to assess the direction of the potential bias. False positives or false negatives are unlikely to have been reported.
- U Indicates the compound or analyte was analyzed for but not detected at or above the stated limit.
- R Data are qualified as rejected. There is a significant potential for the reporting of false negatives or false positives.
- UJ Indicates the compound or analyte was analyzed for but not detected. The sample detection limit is an estimated value.
- A Indicates the finding is based upon technical validation criteria.
- P Indicates the finding is related to a protocol/contractual deviation.
- None Indicates the data was not significantly impacted by the finding, therefore qualification was not required.

## I. Technical Holding Times

All technical holding time requirements were met.

The chain-of-custodies were reviewed for documentation of cooler temperatures. All cooler temperatures met validation criteria.

## II. ICPMS Tune

ICP-MS was not utilized in this SDG.

## III. Calibration

An initial calibration was performed.

The frequency and analysis criteria of the initial calibration verification (ICV) and continuing calibration verification (CCV) were met.

Calibration data were not reviewed for Stage 2A.

## IV. Blanks

Method blanks were reviewed for each matrix as applicable. No chromium was found in the initial, continuing and preparation blanks with the following exceptions:

Method Blank ID	Analyte	Maximum Concentration	Associated Samples
ICB/CCB	Chromium	0.0021 mg/L	ART-7** ART-8** PC-99R2/R3** PC-115R** PC-116R** PC-117** PC-118**

Sample concentrations were compared to concentrations detected in the method blanks as required by the QAPP. No sample data was qualified with the following exceptions:

Sample	Analyte	Reported Concentration	Modified Final Concentration
PC-99R2/R3**	Chromium	0.0031 mg/L	0.02U mg/L
PC-115R**	Chromium	0.0019 mg/L	0.02U mg/L
PC-116R**	Chromium	0.0019 mg/L	0.02U mg/L
PC-117**	Chromium	0.0017 mg/L	0.01U mg/L

Sample	Analyte	Reported Concentration	Modified Final Concentration
PC-118**	Chromium	0.0024 mg/L	0.01U mg/L

Calibration blank data were not reviewed for Stage 2A.

No field blanks were identified in this SDG.

#### V. ICP Interference Check Sample (ICS) Analysis

The frequency of analysis was met.

The criteria for analysis were met.

ICP Interference check sample analysis data were not reviewed for Stage 2A.

#### VI. Matrix Spike Analysis

Matrix spike (MS) and matrix spike duplicate (MSD) samples were reviewed for each matrix as applicable. Percent recoveries (%R) and relative percent differences (RPD) were within QC limits.

#### VII. Duplicate Sample Analysis

Duplicate (DUP) sample analyses were reviewed for each matrix as applicable.

#### VIII. Laboratory Control Samples (LCS)

Laboratory control samples were reviewed for each matrix as applicable. Percent recoveries (%R) and relative percent differences (RPD) were within QC limits.

#### IX. Internal Standards (ICP-MS)

ICP-MS was not utilized in this SDG.

#### X. Furnace Atomic Absorption QC

Graphite furnace atomic absorption was not utilized in this SDG.

#### XI. ICP Serial Dilution

ICP serial dilution was not performed for this SDG.

#### XII. Sample Result Verification

All sample result verifications were acceptable for samples on which a Stage 4 review was performed. Raw data were not evaluated for the samples reviewed by Stage 2A criteria.

### **XIII. Overall Assessment of Data**

Data flags are summarized at the end of this report if data has been qualified.

### **XIV. Field Duplicates**

No field duplicates were identified in this SDG.

**2010 Annual Remedial Performance Sampling  
Chromium - Data Qualification Summary - SDG 340161**

No Sample Data Qualified in this SDG

**2010 Annual Remedial Performance Sampling  
Chromium - Laboratory Blank Data Qualification Summary - SDG 340161**

SDG	Sample	Analyte	Modified Final Concentration	A or P
340161	PC-99R2/R3**	Chromium	0.02U mg/L	A
340161	PC-115R**	Chromium	0.02U mg/L	A
340161	PC-116R**	Chromium	0.02U mg/L	A
340161	PC-117**	Chromium	0.01U mg/L	A
340161	PC-118**	Chromium	0.01U mg/L	A

**2010 Annual Remedial Performance Sampling  
Chromium - Field Blank Data Qualification Summary - SDG 340161**

No Sample Data Qualified in this SDG

Tronox Northgate Henderson

VALIDATION COMPLETENESS WORKSHEET

LDC #: 24203D4

SDG #: 340161

Laboratory: MWH Laboratories

Stage 2B/4

A  
9mg

Date: 12-1-10

Page: 1 of 1

Reviewer: MG

2nd Reviewer: ✓

METHOD: Chromium (EPA SW 846 Method 6010)

The samples listed below were reviewed for each of the following validation areas. Validation findings are noted in attached validation findings worksheets.

	Validation Area		Comments
I.	Technical holding times	A	Sampling dates: 8-4-10
II.	ICP/MS Tune	N	not utilized
III.	Calibration	A	
IV.	Blanks	SW	
V.	ICP Interference Check Sample (ICS) Analysis	A	
VI.	Matrix Spike Analysis	A	MS/MSD (SDG: 340066)
VII.	Duplicate Sample Analysis	N	
VIII.	Laboratory Control Samples (LCS)	A	LCS/LCSD
IX.	Internal Standard (ICP-MS)	N	not utilized
X.	Furnace Atomic Absorption QC	N	" "
XI.	ICP Serial Dilution	N	not performed
XII.	Sample Result Verification	A	Not reviewed for Stage 2B validation.
XIII.	Overall Assessment of Data	A	
XIV.	Field Duplicates	N	
XV.	Field Blanks	N	

Note: A = Acceptable  
 N = Not provided/applicable  
 SW = See worksheet  
 ND = No compounds detected  
 R = Rinsate  
 FB = Field blank  
 D = Duplicate  
 TB = Trip blank  
 EB = Equipment blank

Validated Samples: \*\* Indicates sample underwent Stage 4 validation  
 all water

1	ART-1	11	PC-117**	21	PC-118MSD	31	
2	ART-2	12	PC-118**	22	PBW1	32	
3	ART-3	13	PC-119	23	PBW2	33	
4	ART-4	14	PC-120	24		34	
5	ART-6	15	PC-121	25		35	
6	ART-7**	16	PC-133	26		36	
7	ART-8**	17	ART-9	27		37	
8	PC-99R2/R3**	18	PC-117MS	28		38	
9	PC-115R**	19	PC-117MSD	29		39	
10	PC-116R**	20	PC-118MS	30		40	

Notes: \_\_\_\_\_  
 \_\_\_\_\_  
 \_\_\_\_\_



**Method:Metals (EPA SW 846 Method 6010B/7000/6020)**

Validation Area	Yes	No	NA	Findings/Comments
<b>I. Technical holding times</b>				
All technical holding times were met.	✓			
Cooler temperature criteria was met.	✓			
<b>II. ICP/MS Tune</b>				
Were all isotopes in the tuning solution mass resolution within 0.1 amu?			✓	
Were %RSD of isotopes in the tuning solution $\leq 5\%$ ?			✓	
<b>III. Calibration</b>				
Were all instruments calibrated daily, each set-up time?	✓			
Were the proper number of standards used?	✓			
Were all initial and continuing calibration verification %Rs within the 90-110% (80-120% for mercury) QC limits?	✓			
Were all initial calibration correlation coefficients $> 0.995$ ?	✓			
<b>IV. Blanks</b>				
Was a method blank associated with every sample in this SDG?	✓			
Was there contamination in the method blanks? If yes, please see the Blanks validation completeness worksheet.	✓			
<b>V. ICP Interference Check Sample</b>				
Were ICP interference check samples performed daily?	✓			
Were the AB solution percent recoveries (%R) with the 80-120% QC limits?	✓			
<b>VI. Matrix spike/Matrix spike duplicates</b>				
Were a matrix spike (MS) and duplicate (DUP) analyzed for each matrix in this SDG? If no, indicate which matrix does not have an associated MS/MSD or MS/DUP. Soil / Water.	✓			
Were the MS/MSD percent recoveries (%R) and the relative percent differences (RPD) within the 75-125 QC limits? If the sample concentration exceeded the spike concentration by a factor of 4 or more, no action was taken.	✓			
Were the MS/MSD or duplicate relative percent differences (RPD) $\leq 20\%$ for waters and $\leq 35\%$ for soil samples? A control limit of $\pm RL$ ( $\pm 2X RL$ for soil) was used for samples that were $\leq 5X$ the RL, including when only one of the duplicate sample values were $< 5X$ the RL.	✓			
<b>VII. Laboratory control samples</b>				
Was an LCS analyzed for this SDG?	✓			
Was an LCS analyzed per extraction batch?	✓			
Were the LCS percent recoveries (%R) and relative percent difference (RPD) within the 80-120% QC limits for water samples and laboratory established QC limits for soils?	✓			

Validation Area	Yes	No	NA	Findings/Comments
<b>VIII. Furnace Atomic Absorption QC</b>				
If MSA was performed, was the correlation coefficients > 0.995?			✓	
Do all applicable analyses have duplicate injections? (Level IV only)			✓	
For sample concentrations > RL, are applicable duplicate injection RSD values < 20%? (Level IV only)			✓	
Were analytical spike recoveries within the 85-115% QC limits?			✓	
<b>IX. ICP Serial Dilution</b>				
Was an ICP serial dilution analyzed if analyte concentrations were > 50X the MDL (ICP)/>100X the MDL(ICP/MS)?		✓		
Were all percent differences (%Ds) < 10%?			✓	
Was there evidence of negative interference? If yes, professional judgement will be used to qualify the data.			✓	
<b>X. Internal Standards (EPA SW 846 Method 6020/EPA 200.8)</b>				
Were all the percent recoveries (%R) within the 30-120% (6020)/60-125% (200.8) of the intensity of the internal standard in the associated initial calibration?			✓	
If the %Rs were outside the criteria, was a reanalysis performed?			✓	
<b>XI. Regional Quality Assurance and Quality Control</b>				
Were performance evaluation (PE) samples performed?		✓		
Were the performance evaluation (PE) samples within the acceptance limits?			✓	
<b>XII. Sample Result Verification</b>				
Were RLs adjusted to reflect all sample dilutions and dry weight factors applicable to level IV validation?	✓			
<b>XIII. Overall assessment of data</b>				
Overall assessment of data was found to be acceptable.	✓			
<b>XIV. Field duplicates</b>				
Field duplicate pairs were identified in this SDG.		✓		
Target analytes were detected in the field duplicates.			✓	
<b>XV. Field blanks</b>				
Field blanks were identified in this SDG.		✓		
Target analytes were detected in the field blanks.			✓	

VALIDATION FINDINGS WORKSHEET  
 PB/ICB/CCB QUALIFIED SAMPLES

METHOD: Trace metals (EPA SW 864 Method 6010B/6020/7000)

Sample Concentration units, unless otherwise noted: mg/L

Soil preparation factor applied: NA

Associated Samples: 6 → 12

Code: bl

Analyte	Maximum PB <sup>a</sup> (mg/Kg)	Maximum PB <sup>a</sup> (mg/L)	Maximum ICB/CCB <sup>a</sup> (mg/L)	Action Limit	1	8	9	10	11	12	13	14	15
Cr			0.0021		0.0048/ 0.02U	0.0031/ 0.02U	0.0019/ 0.02U	0.0019/ 0.02U	0.0017/ 0.01U	0.0024/ 0.01U	0.0014/ 0.02U	0.0016/ 0.02U	0.0018/ 0.02U

Samples with analyte concentrations within five times the associated ICB, CCB or PB concentration are listed above with the identifications from the Validation Completeness Worksheet. These sample results were qualified as not detected, "U".

Note: a - The listed analyte concentration is the highest ICB, CCB, or PB detected in the analysis of each element.

*SM*

*SM*

LDC #: 24203D4

**VALIDATION FINDINGS WORKSHEET**  
**Initial and Continuing Calibration Calculation Verification**

Page: 1 of 1  
 Reviewer: MG  
 2nd Reviewer: W

**METHOD:** Trace Metals (EPA SW 846 Method 6010/6020/7000)

An initial and continuing calibration verification percent recovery (%R) was recalculated for each type of analysis using the following formula:

$\%R = \frac{\text{Found}}{\text{True}} \times 100$  Where, Found = concentration (in ug/L) of each analyte measured in the analysis of the ICV or CCV solution  
 True = concentration (in ug/L) of each analyte in the ICV or CCV source

Standard ID	Type of Analysis	Element	Found (ug/L)	True (ug/L)	Recalculated		Reported		Acceptable (Y/N)
					%R	%R	%R	%R	
1933 ICV	ICP (Initial calibration)	Cr	10033	10000	100	100	100	100	Y
	ICP/MS (Initial calibration)								
	CVAA (Initial calibration)								
1951 CCV	ICP (Continuing calibration)	Cr	4933.9	5000	98.7	98.6	98.6	98.6	↓
	ICP/MS (Continuing calibration)								
	CVAA (Continuing calibration)								
	GFAA (Initial calibration)								
	GFAA (Continuing calibration)								

Comments: Refer to Calibration Verification findings worksheet for list of qualifications and associated samples when reported results do not agree within 10.0% of the recalculated results.

**VALIDATION FINDINGS WORKSHEET**  
**Level IV Recalculation Worksheet**

**METHOD:** Trace Metals (EPA SW 846 Method 6010/6020/7000)

Percent recoveries (%R) for an ICP interference check sample, a laboratory control sample and a matrix spike sample were recalculated using the following formula:

$$\%R = \frac{\text{Found} \times 100}{\text{True}}$$

Where, Found = Concentration of each analyte measured in the analysis of the sample. For the matrix spike calculation,  
 Found = SSR (spiked sample result) - SR (sample result).  
 True = Concentration of each analyte in the source.

A sample and duplicate relative percent difference (RPD) was recalculated using the following formula:

$$RPD = \frac{|S-D|}{(S+D)/2} \times 100$$

Where, S = Original sample concentration  
 D = Duplicate sample concentration

An ICP serial dilution percent difference (%D) was recalculated using the following formula:

$$\%D = \frac{|I-SDR|}{I} \times 100$$

Where, I = Initial Sample Result (mg/L)  
 SDR = Serial Dilution Result (mg/L) (Instrument Reading x 5)

Sample ID	Type of Analysis	Element	Found / S / I (units)	True / D / SDR (units)	Recalculated		Acceptable (Y/N)
					%R / RPD / %D	Reported %R / RPD / %D	
1943 ICSA B	ICP interference check	Cr	0.24070 mg/L	0.25 (mg/L)	96.3	96.2	Y
0118 LCS 2	Laboratory control sample	Cr	1.0055 (mg/L)	1.0 (mg/L)	101	101	Y
0128 18	Matrix spike	Cr	0.97855 (mg/L) (SSR-SR)	1.0 (mg/L)	98	98	Y
0128 / 0132 18 / 19	Duplicate	Cr	0.98025 (mg/L)	1.0171 (mg/L)	3.7	4.1	Y
—	ICP serial dilution	—	—	—	—	—	—

Comments: Refer to appropriate worksheet for list of qualifications and associated samples when reported results do not agree within 10.0% of the recalculated results.



## Laboratory Data Consultants, Inc. Data Validation Report

**Project/Site Name:** 2010 Annual Remedial Performance Sampling  
**Collection Date:** August 9 through August 12, 2010  
**LDC Report Date:** December 19, 2010  
**Matrix:** Water  
**Parameters:** Chromium  
**Validation Level:** Stage 2A  
**Laboratory:** MWH Laboratories, Inc.  
**Sample Delivery Group (SDG):** 340887

### Sample Identification

M-83	MW-K5
M-87	PC-91
PC-86	PC-97
PC-90	PC-18
PC-56	PC-55
PC-58	PC-101R
PC-59	ART-7B
PC-60	PC-92
PC-62	PC-94
PC-68	M-83MS
PC-122	M-83MSD
MW-K4	PC-86MS
ARP-1	PC-86MSD
ARP-2A	MW-K5MS
ARP-3A	MW-K5MSD
ARP-4A	
ARP-5A	
ARP-6B	
ARP-7	
PC-53	

## Introduction

This data review covers 35 water samples listed on the cover sheet including dilutions and reanalysis as applicable. The analyses were per EPA SW 846 Method 6010 for Chromium.

This review follows a modified outline of the USEPA Contract Laboratory Program National Functional Guidelines for Inorganic Data Review (October 2004) and the EPA Region 9 Superfund Data Evaluation/Validation Guidance, NDEP guidance (May 2006).

A qualification summary table is provided at the end of this report if data has been qualified. Flags are classified as P (protocol) or A (advisory) to indicate whether the flag is due to a laboratory deviation from a specified protocol or is of technical advisory nature.

Raw data were not reviewed for this SDG. The review was based on QC data.

The following are definitions of the data qualifiers:

- J+ Data are qualified as estimated, with a high bias likely to occur. False positives or false negatives are unlikely to have been reported.
- J- Data are qualified as estimated, with a low bias likely to occur. False positives or false negatives are unlikely to have been reported.
- J Data are qualified as estimated; it is not possible to assess the direction of the potential bias. False positives or false negatives are unlikely to have been reported.
- U Indicates the compound or analyte was analyzed for but not detected at or above the stated limit.
- R Data are qualified as rejected. There is a significant potential for the reporting of false negatives or false positives.
- UJ Indicates the compound or analyte was analyzed for but not detected. The sample detection limit is an estimated value.
- A Indicates the finding is based upon technical validation criteria.
- P Indicates the finding is related to a protocol/contractual deviation.
- None Indicates the data was not significantly impacted by the finding, therefore qualification was not required.



## **I. Technical Holding Times**

All technical holding time requirements were met.

The chain-of-custodies were reviewed for documentation of cooler temperatures. All cooler temperatures met validation criteria.

## **II. ICPMS Tune**

ICP-MS was not utilized in this SDG.

## **III. Calibration**

Calibration data were not reviewed for Stage 2A.

## **IV. Blanks**

Method blanks were reviewed for each matrix as applicable. No chromium was found in the preparation blanks.

No field blanks were identified in this SDG.

## **V. ICP Interference Check Sample (ICS) Analysis**

ICP Interference check sample analysis data were not reviewed for Stage 2A.

## **VI. Matrix Spike Analysis**

Matrix spike (MS) and matrix spike duplicate (MSD) samples were reviewed for each matrix as applicable. Percent recoveries (%R) and relative percent differences (RPD) were within QC limits.

## **VII. Duplicate Sample Analysis**

Duplicate (DUP) sample analyses were reviewed for each matrix as applicable.

## **VIII. Laboratory Control Samples (LCS)**

Laboratory control samples were reviewed for each matrix as applicable. Percent recoveries (%R) and relative percent differences (RPD) were within QC limits.

## **IX. Internal Standards (ICP-MS)**

ICP-MS was not utilized in this SDG.

#### **X. Furnace Atomic Absorption QC**

Graphite furnace atomic absorption was not utilized in this SDG.

#### **XI. ICP Serial Dilution**

ICP serial dilution analysis data were not reviewed for Stage 2A.

#### **XII. Sample Result Verification**

Raw data were not reviewed for this SDG.

#### **XIII. Overall Assessment of Data**

Data flags are summarized at the end of this report if data has been qualified.

#### **XIV. Field Duplicates**

No field duplicates were identified in this SDG.

**2010 Annual Remedial Performance Sampling  
Chromium - Data Qualification Summary - SDG 340887**

No Sample Data Qualified in this SDG

**2010 Annual Remedial Performance Sampling  
Chromium - Laboratory Blank Data Qualification Summary - SDG 340887**

No Sample Data Qualified in this SDG

**2010 Annual Remedial Performance Sampling  
Chromium - Field Blank Data Qualification Summary - SDG 340887**

No Sample Data Qualified in this SDG

Tronox Northgate Henderson

VALIDATION COMPLETENESS WORKSHEET

LDC #: 24203J4

SDG #: 340887

Laboratory: MWH Laboratories

Stage 2BA

gmf

Date: 12-2-10

Page: 1 of 1

Reviewer: MG

2nd Reviewer: [Signature]

METHOD: Chromium (EPA SW 846 Method 6010)

The samples listed below were reviewed for each of the following validation areas. Validation findings are noted in attached validation findings worksheets.

	Validation Area		Comments
I.	Technical holding times	A	Sampling dates: 8-9-10 through 8-12-10
II.	ICP/MS Tune	N	not utilized
III.	Calibration	N	
IV.	Blanks	A	
V.	ICP Interference Check Sample (ICS) Analysis	N	
VI.	Matrix Spike Analysis	A	MS/MSD
VII.	Duplicate Sample Analysis	N	
VIII.	Laboratory Control Samples (LCS)	A	LCS/LCSD
IX.	Internal Standard (ICP-MS)	N	not utilized
X.	Furnace Atomic Absorption QC	N	" "
XI.	ICP Serial Dilution	N	not performed
XII.	Sample Result Verification	N	
XIII.	Overall Assessment of Data	A	
XIV.	Field Duplicates	N	
XV.	Field Blanks	N	

Note: A = Acceptable  
N = Not provided/applicable  
SW = See worksheet

ND = No compounds detected  
R = Rinsate  
FB = Field blank

D = Duplicate  
TB = Trip blank  
EB = Equipment blank

Validated Samples:  
all water

1	M-83	11	PC-122	21	MW-K5	31	M-83MSD
2	M-87	12	MW-K4	22	PC-91	32	PC-86MS
3	PC-86	13	ARP-1	23	PC-97	33	PC-86MSD
4	PC-90	14	ARP-2A	24	PC-18	34	MW-K5MS
5	PC-56	15	ARP-3A	25	PC-55	35	MW-K5MSD
6	PC-58	16	ARP-4A	26	PC-101R	36	
7	PC-59	17	ARP-5A	27	ART-7B	37	
8	PC-60	18	ARP-6B	28	PC-92	38	PBW1
9	PC-62	19	ARP-7	29	PC-94	39	PBW2
10	PC-68	20	PC-53	30	M-83MS	40	PBW3

Notes: \_\_\_\_\_  
\_\_\_\_\_  
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**Laboratory Data Consultants, Inc.  
Data Validation Report**

**Project/Site Name:** 2010 Annual Remedial Performance Sampling  
**Collection Date:** August 23, 2010  
**LDC Report Date:** December 19, 2010  
**Matrix:** Water  
**Parameters:** Chromium  
**Validation Level:** Stage 2A  
**Laboratory:** MWH Laboratories, Inc.  
**Sample Delivery Group (SDG):** 341684

**Sample Identification**

PC-144  
PC-145

## Introduction

This data review covers 2 water samples listed on the cover sheet including dilutions and reanalysis as applicable. The analyses were per EPA SW 846 Method 6010 for Chromium.

This review follows a modified outline of the USEPA Contract Laboratory Program National Functional Guidelines for Inorganic Data Review (October 2004) and the EPA Region 9 Superfund Data Evaluation/Validation Guidance, NDEP guidance (May 2006).

A qualification summary table is provided at the end of this report if data has been qualified. Flags are classified as P (protocol) or A (advisory) to indicate whether the flag is due to a laboratory deviation from a specified protocol or is of technical advisory nature.

Raw data were not reviewed for this SDG. The review was based on QC data.

The following are definitions of the data qualifiers:

J+ Data are qualified as estimated, with a high bias likely to occur. False positives or false negatives are unlikely to have been reported.

J- Data are qualified as estimated, with a low bias likely to occur. False positives or false negatives are unlikely to have been reported.

J Data are qualified as estimated; it is not possible to assess the direction of the potential bias. False positives or false negatives are unlikely to have been reported.

U Indicates the compound or analyte was analyzed for but not detected at or above the stated limit.

R Data are qualified as rejected. There is a significant potential for the reporting of false negatives or false positives.

UJ Indicates the compound or analyte was analyzed for but not detected. The sample detection limit is an estimated value.

A Indicates the finding is based upon technical validation criteria.

P Indicates the finding is related to a protocol/contractual deviation.

None Indicates the data was not significantly impacted by the finding, therefore qualification was not required.

## **I. Technical Holding Times**

All technical holding time requirements were met.

The chain-of-custodies were reviewed for documentation of cooler temperatures. All cooler temperatures met validation criteria.

## **II. ICPMS Tune**

ICP-MS was not utilized in this SDG.

## **III. Calibration**

Calibration data were not reviewed for Stage 2A.

## **IV. Blanks**

Method blanks were reviewed for each matrix as applicable. No chromium was found in the preparation blanks.

No field blanks were identified in this SDG.

## **V. ICP Interference Check Sample (ICS) Analysis**

ICP Interference check sample analysis data were not reviewed for Stage 2A.

## **VI. Matrix Spike Analysis**

The laboratory has indicated that there were no matrix spike (MS) and matrix spike duplicate (MSD) analyses specified for the samples in this SDG, and therefore matrix spike and matrix spike duplicate analyses were not performed for this SDG.

## **VII. Duplicate Sample Analysis**

The laboratory has indicated that there were no duplicate (DUP) analyses specified for the samples in this SDG, and therefore duplicate analyses were not performed for this SDG.

## **VIII. Laboratory Control Samples (LCS)**

Laboratory control samples were reviewed for each matrix as applicable. Percent recoveries (%R) and relative percent differences (RPD) were within QC limits.

## **IX. Internal Standards (ICP-MS)**

ICP-MS was not utilized in this SDG.

#### **X. Furnace Atomic Absorption QC**

Graphite furnace atomic absorption was not utilized in this SDG.

#### **XI. ICP Serial Dilution**

ICP serial dilution analysis data were not reviewed for Stage 2A.

#### **XII. Sample Result Verification**

Raw data were not reviewed for this SDG.

#### **XIII. Overall Assessment of Data**

Data flags are summarized at the end of this report if data has been qualified.

#### **XIV. Field Duplicates**

No field duplicates were identified in this SDG.



**2010 Annual Remedial Performance Sampling  
Chromium - Data Qualification Summary - SDG 341684**

No Sample Data Qualified in this SDG

**2010 Annual Remedial Performance Sampling  
Chromium - Laboratory Blank Data Qualification Summary - SDG 341684**

No Sample Data Qualified in this SDG

**2010 Annual Remedial Performance Sampling  
Chromium - Field Blank Data Qualification Summary - SDG 341684**

No Sample Data Qualified in this SDG

Tronox Northgate Henderson

LDC #: 24203K4

VALIDATION COMPLETENESS WORKSHEET

Date: 12-2-10

SDG #: 341684

Stage 2B A

Page: 1 of 1

Laboratory: MWH Laboratories

MM

Reviewer: MG

2nd Reviewer: [Signature]

METHOD: Chromium (EPA SW 846 Method 6010)

The samples listed below were reviewed for each of the following validation areas. Validation findings are noted in attached validation findings worksheets.

	Validation Area		Comments
I.	Technical holding times	A	Sampling dates: 8-23-10
II.	ICP/MS Tune	N	not utilized
III.	Calibration	N	
IV.	Blanks	A	
V.	ICP Interference Check Sample (ICS) Analysis	N	
VI.	Matrix Spike Analysis	N	client specified
VII.	Duplicate Sample Analysis	N	" "
VIII.	Laboratory Control Samples (LCS)	A	LCS/LCSD
IX.	Internal Standard (ICP-MS)	N	not utilized
X.	Furnace Atomic Absorption QC	N	" "
XI.	ICP Serial Dilution	N	not performed
XII.	Sample Result Verification	N	
XIII.	Overall Assessment of Data	A	
XIV.	Field Duplicates	N	
XV.	Field Blanks	N	

Note: A = Acceptable  
N = Not provided/applicable  
SW = See worksheet

ND = No compounds detected  
R = Rinsate  
FB = Field blank

D = Duplicate  
TB = Trip blank  
EB = Equipment blank

Validated Samples:  
all water

1	PC-144	11		21		31	
2	PC-145	12		22		32	
3	PBW	13		23		33	
4		14		24		34	
5		15		25		35	
6		16		26		36	
7		17		27		37	
8		18		28		38	
9		19		29		39	
10		20		30		40	

Notes: \_\_\_\_\_  
\_\_\_\_\_  
\_\_\_\_\_

## Laboratory Data Consultants, Inc. Data Validation Report

**Project/Site Name:** 2010 Annual Remedial Performance Sampling  
**Collection Date:** August 1 through August 2, 2010  
**LDC Report Date:** December 19, 2010  
**Matrix:** Water  
**Parameters:** Chromium  
**Validation Level:** Stage 2A  
**Laboratory:** MWH Laboratories, Inc.  
**Sample Delivery Group (SDG):** 339791

### Sample Identification

I-O	PC-132
I-P	M-96
I-H	PC-54
I-U	PC-37
I-T	PC-71
I-G	PC-72
I-F	PC-73
I-N	M-23
I-E	VD080210
I-M	FB080210V
I-D	PC-71MS
PC-123	PC-71MSD
PC-124	PC-73MS
PC-125	PC-73MSD
PC-126	M-23MS
PC-127	M-23MSD
PC-128	VD080210MS
PC-129	VD080210MSD
PC-130	
PC-131	

## Introduction

This data review covers 38 water samples listed on the cover sheet including dilutions and reanalysis as applicable. The analyses were per EPA SW 846 Method 6010 for Chromium.

This review follows a modified outline of the USEPA Contract Laboratory Program National Functional Guidelines for Inorganic Data Review (October 2004) and the EPA Region 9 Superfund Data Evaluation/Validation Guidance, NDEP guidance (May 2006).

A qualification summary table is provided at the end of this report if data has been qualified. Flags are classified as P (protocol) or A (advisory) to indicate whether the flag is due to a laboratory deviation from a specified protocol or is of technical advisory nature.

Raw data were not reviewed for this SDG. The review was based on QC data.

The following are definitions of the data qualifiers:

- J+ Data are qualified as estimated, with a high bias likely to occur. False positives or false negatives are unlikely to have been reported.
- J- Data are qualified as estimated, with a low bias likely to occur. False positives or false negatives are unlikely to have been reported.
- J Data are qualified as estimated; it is not possible to assess the direction of the potential bias. False positives or false negatives are unlikely to have been reported.
- U Indicates the compound or analyte was analyzed for but not detected at or above the stated limit.
- R Data are qualified as rejected. There is a significant potential for the reporting of false negatives or false positives.
- UJ Indicates the compound or analyte was analyzed for but not detected. The sample detection limit is an estimated value.
- A Indicates the finding is based upon technical validation criteria.
- P Indicates the finding is related to a protocol/contractual deviation.
- None Indicates the data was not significantly impacted by the finding, therefore qualification was not required.

## I. Technical Holding Times

All technical holding time requirements were met.

The chain-of-custodies were reviewed for documentation of cooler temperatures. All cooler temperatures met validation criteria.

## II. ICPMS Tune

ICP-MS was not utilized in this SDG.

## III. Calibration

Calibration data were not reviewed for Stage 2A.

## IV. Blanks

Method blanks were reviewed for each matrix as applicable. No chromium was found in the preparation blanks.

Sample FB080210V was identified as a field blank. No chromium was found in this blank with the following exceptions:

Field Blank ID	Sampling Date	Analyte	Concentration	Associated Samples
FB080210V	8/2/10	Chromium	0.006 mg/L	I-O I-P I-H I-U I-T I-G I-F I-N I-E I-M I-D PC-54 PC-37 PC-71 PC-72 PC-73 M-23 VD080210

Sample concentrations were compared to concentrations detected in the field blanks as required by the QAPP. No sample data was qualified.

## V. ICP Interference Check Sample (ICS) Analysis

ICP Interference check sample analysis data were not reviewed for Stage 2A.

## VI. Matrix Spike Analysis

Matrix spike (MS) and matrix spike duplicate (MSD) samples were reviewed for each matrix as applicable. Percent recoveries (%R) and relative percent differences (RPD) were within QC limits.

## VII. Duplicate Sample Analysis

Duplicate (DUP) sample analyses were reviewed for each matrix as applicable.

## VIII. Laboratory Control Samples (LCS)

Laboratory control samples were reviewed for each matrix as applicable. Percent recoveries (%R) and relative percent differences (RPD) were within QC limits.

## IX. Internal Standards (ICP-MS)

ICP-MS was not utilized in this SDG.

## X. Furnace Atomic Absorption QC

Graphite furnace atomic absorption was not utilized in this SDG.

## XI. ICP Serial Dilution

ICP serial dilution analysis data were not reviewed for Stage 2A.

## XII. Sample Result Verification

Raw data were not reviewed for this SDG.

## XIII. Overall Assessment of Data

Data flags are summarized at the end of this report if data has been qualified.

## XIV. Field Duplicates

Samples M-23 and VD080210 were identified as field duplicates. No contaminant concentrations were detected in any of the samples with the following exceptions:

Analyte	Concentration (mg/L)		RPD (Limits)	Difference (Limits)	Flags	A or P
	M-23	VD080210				
Chromium	0.54	0.51	6 (≤30)	-	-	-

**2010 Annual Remedial Performance Sampling  
Chromium - Data Qualification Summary - SDG 339791**

No Sample Data Qualified in this SDG

**2010 Annual Remedial Performance Sampling  
Chromium - Laboratory Blank Data Qualification Summary - SDG 339791**

No Sample Data Qualified in this SDG

**2010 Annual Remedial Performance Sampling  
Chromium - Field Blank Data Qualification Summary - SDG 339791**

No Sample Data Qualified in this SDG

Tronox Northgate Henderson

VALIDATION COMPLETENESS WORKSHEET

LDC #: 24203N4

SDG #: 339791

Laboratory: MWH Laboratories

Stage 2 ~~B~~ A

MH

Date: 12-15-10

Page: 1 of 1

Reviewer: MG

2nd Reviewer: W

METHOD: Chromium (EPA SW 846 Method 6010)

The samples listed below were reviewed for each of the following validation areas. Validation findings are noted in attached validation findings worksheets.

	Validation Area		Comments
I.	Technical holding times	A	Sampling dates: 8-1-10 through 8-2-10
II.	ICP/MS Tune	N	not utilized
III.	Calibration	N	
IV.	Blanks	A	
V.	ICP Interference Check Sample (ICS) Analysis	N	
VI.	Matrix Spike Analysis	A	MS/MSD
VII.	Duplicate Sample Analysis	N	
VIII.	Laboratory Control Samples (LCS)	A	LCS/LCSD
IX.	Internal Standard (ICP-MS)	N	not utilized
X.	Furnace Atomic Absorption QC	N	" "
XI.	ICP Serial Dilution	N	not performed
XII.	Sample Result Verification	N	
XIII.	Overall Assessment of Data	A	
XIV.	Field Duplicates	SW	D = 28 + 29
XV.	Field Blanks	SW	FB = 30

Note: A = Acceptable  
N = Not provided/applicable  
SW = See worksheet

ND = No compounds detected  
R = Rinsate  
FB = Field blank

D = Duplicate  
TB = Trip blank  
EB = Equipment blank

Validated Samples:  
all water

1	1	I-O	11	2	I-D	21	2	PC-132	31	1	PC-71MS
2	2	I-P	12	1	PC-123	22	1	M-96	32	1	PC-71MSD
3	1	I-H	13	2	PC-124	23	2	PC-54	33	2	PC-73MS
4	1	I-U	14	2	PC-125	24	2	PC-37	34	2	PC-73MSD
5	2	I-T	15	1	PC-126	25	1	PC-71	35	2	M-23MS
6	1	I-G	16	1	PC-127	26	2	PC-72	36	2	M-23MSD
7	1	I-F	17	1	PC-128	27	2	PC-73	37	1	VD080210MS
8	1	I-N	18	1	PC-129	28	2	M-23	38	1	VD080210MSD
9	1	I-E	19	1	PC-130	29	1	VD080210	39	1	PBW1
10	1	I-M	20	1	PC-131	30	1	FB080210V	40	2	PBW2

Notes:

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LDC#: 24203N4

**VALIDATION FINDINGS WORKSHEET**  
**Field Duplicates**

Page: 1 of 1  
Reviewer: MG  
2nd Reviewer: [Signature]

**METHOD:** Metals (EPA Method 6010B/6020/7000)

- Y N NA Were field duplicate pairs identified in this SDG?  
 Y N NA Were target analytes detected in the field duplicate pairs?

Analyte	Concentration (mg/L)		( $\leq 30$ )	(mg/L)	(mg/L)	Qualifications (Parent Only)
	28	29	RPD	Difference	Limits	
Chromium	0.54	0.51	6			

V:\FIELD DUPLICATES\FD\_inorganic\24203N4.wpd

## Laboratory Data Consultants, Inc. Data Validation Report

**Project/Site Name:** 2010 Annual Remedial Performance Sampling  
**Collection Date:** August 2 through August 3, 2010  
**LDC Report Date:** December 19, 2010  
**Matrix:** Water  
**Parameters:** Chromium  
**Validation Level:** Stage 2A  
**Laboratory:** MWH Laboratories, Inc.  
**Sample Delivery Group (SDG):** 339977

### Sample Identification

I-Q  
I-C  
I-S  
I-L  
I-R  
I-B  
I-AR  
I-AB  
I-AA  
M-131  
M-57A  
M-79  
M-69  
M-135  
VD-080310  
EB080310V  
I-ABMS  
I-ABMSD  
M-135MS  
M-135MSD

## Introduction

This data review covers 20 water samples listed on the cover sheet including dilutions and reanalysis as applicable. The analyses were per EPA SW 846 Method 6010 for Chromium.

This review follows a modified outline of the USEPA Contract Laboratory Program National Functional Guidelines for Inorganic Data Review (October 2004) and the EPA Region 9 Superfund Data Evaluation/Validation Guidance, NDEP guidance (May 2006).

A qualification summary table is provided at the end of this report if data has been qualified. Flags are classified as P (protocol) or A (advisory) to indicate whether the flag is due to a laboratory deviation from a specified protocol or is of technical advisory nature.

Raw data were not reviewed for this SDG. The review was based on QC data.

The following are definitions of the data qualifiers:

- J+ Data are qualified as estimated, with a high bias likely to occur. False positives or false negatives are unlikely to have been reported.
- J- Data are qualified as estimated, with a low bias likely to occur. False positives or false negatives are unlikely to have been reported.
- J Data are qualified as estimated; it is not possible to assess the direction of the potential bias. False positives or false negatives are unlikely to have been reported.
- U Indicates the compound or analyte was analyzed for but not detected at or above the stated limit.
- R Data are qualified as rejected. There is a significant potential for the reporting of false negatives or false positives.
- UJ Indicates the compound or analyte was analyzed for but not detected. The sample detection limit is an estimated value.
- A Indicates the finding is based upon technical validation criteria.
- P Indicates the finding is related to a protocol/contractual deviation.
- None Indicates the data was not significantly impacted by the finding, therefore qualification was not required.

## I. Technical Holding Times

All technical holding time requirements were met.

The chain-of-custodies were reviewed for documentation of cooler temperatures. All cooler temperatures met validation criteria.

## II. ICPMS Tune

ICP-MS was not utilized in this SDG.

## III. Calibration

Calibration data were not reviewed for Stage 2A.

## IV. Blanks

Method blanks were reviewed for each matrix as applicable. No chromium was found in the preparation blanks.

Sample EB080310V was identified as an equipment blank. No chromium was found in this blank with the following exceptions:

Equipment Blank ID	Sampling Date	Analyte	Concentration	Associated Samples
EB080310V	8/3/10	Chromium	0.0014 mg/L	M-131 M-57A M-79 M-69 M-135 VD-080310

Sample FB080210V (from SDG 339791) was identified as a field blank. No chromium was found in this blank with the following exceptions:

Field Blank ID	Sampling Date	Analyte	Concentration	Associated Samples
FB080210V	8/2/10	Chromium	0.006 mg/L	I-Q I-C I-S I-L I-R I-B I-AR I-AB I-AA

Sample concentrations were compared to concentrations detected in the field blanks as required by the QAPP. No sample data was qualified.

## **V. ICP Interference Check Sample (ICS) Analysis**

ICP Interference check sample analysis data were not reviewed for Stage 2A.

## **VI. Matrix Spike Analysis**

Matrix spike (MS) and matrix spike duplicate (MSD) samples were reviewed for each matrix as applicable. Percent recoveries (%R) and relative percent differences (RPD) were within QC limits.

## **VII. Duplicate Sample Analysis**

Duplicate (DUP) sample analyses were reviewed for each matrix as applicable.

## **VIII. Laboratory Control Samples (LCS)**

Laboratory control samples were reviewed for each matrix as applicable. Percent recoveries (%R) and relative percent differences (RPD) were within QC limits.

## **IX. Internal Standards (ICP-MS)**

ICP-MS was not utilized in this SDG.

## **X. Furnace Atomic Absorption QC**

Graphite furnace atomic absorption was not utilized in this SDG.

## **XI. ICP Serial Dilution**

ICP serial dilution analysis data were not reviewed for Stage 2A.

## **XII. Sample Result Verification**

Raw data were not reviewed for this SDG.

## **XIII. Overall Assessment of Data**

Data flags are summarized at the end of this report if data has been qualified.

## **XIV. Field Duplicates**

Samples M-131 and VD-080310 were identified as field duplicates. No contaminant concentrations were detected in any of the samples with the following exceptions:

Analyte	Concentration (mg/L)		RPD (Limits)	Difference (Limits)	Flags	A or P
	M-131	VD-080310				
Chromium	0.11	0.098	12 (≤30)	-	-	-

**2010 Annual Remedial Performance Sampling  
Chromium - Data Qualification Summary - SDG 339977**

No Sample Data Qualified in this SDG

**2010 Annual Remedial Performance Sampling  
Chromium - Laboratory Blank Data Qualification Summary - SDG 339977**

No Sample Data Qualified in this SDG

**2010 Annual Remedial Performance Sampling  
Chromium - Equipment Blank Data Qualification Summary - SDG 339977**

No Sample Data Qualified in this SDG

**2010 Annual Remedial Performance Sampling  
Chromium - Field Blank Data Qualification Summary - SDG 339977**

No Sample Data Qualified in this SDG



Tronox Northgate Henderson

LDC #: 2420304

VALIDATION COMPLETENESS WORKSHEET

Date: 12-15-10

SDG #: 339977

Stage 2B A

Page: 1 of 1

Laboratory: MWH Laboratories

9mM

Reviewer: MG

2nd Reviewer:

METHOD: Chromium (EPA SW 846 Method 6010)

The samples listed below were reviewed for each of the following validation areas. Validation findings are noted in attached validation findings worksheets.

	Validation Area		Comments
I.	Technical holding times	A	Sampling dates: 8-2-10 through 8-3-10
II.	ICP/MS Tune	N	not utilized
III.	Calibration	N	
IV.	Blanks	A	
V.	ICP Interference Check Sample (ICS) Analysis	N	
VI.	Matrix Spike Analysis	A	MS/MSD (SDG: 339791)
VII.	Duplicate Sample Analysis	N	
VIII.	Laboratory Control Samples (LCS)	A	LCS/LCSD
IX.	Internal Standard (ICP-MS)	N	not utilized
X.	Furnace Atomic Absorption QC	N	" "
XI.	ICP Serial Dilution	N	not performed
XII.	Sample Result Verification	N	
XIII.	Overall Assessment of Data	A	
XIV.	Field Duplicates	SW	D = 10 + 15
XV.	Field Blanks	SW	EB = 16, FB = FB080210V (SDG: 339791)

Note: A = Acceptable  
N = Not provided/applicable  
SW = See worksheet

ND = No compounds detected  
R = Rinsate  
FB = Field blank

D = Duplicate  
TB = Trip blank  
EB = Equipment blank

Validated Samples:

all water

1	1	I-Q	11	1	M-57A	21	1	PBW1	31	
2	2	I-C	12	1	M-79	22	2	PBW2	32	
3	2	I-S	13	1	M-69	23	3	PBW3	33	
4	1	I-L	14	2	M-135	24			34	
5	2	I-R	15	1	VD-080310	25			35	
6	3	I-B	16	2	EB080310V	26			36	
7	1	I-AR	17	2	I-ABMS	27			37	
8	2	I-AB	18	2	I-ABMSD	28			38	
9	1	I-AA	19	2	M-135MS	29			39	
10	1	M-131	20	2	M-135MSD	30			40	

Notes: \_\_\_\_\_  
 \_\_\_\_\_  
 \_\_\_\_\_

**VALIDATION FINDINGS WORKSHEET**  
**Field Blanks**

**METHOD:** Trace Metals (EPA SW846 6010B/7000)  
 Y/N N/A Were field blanks identified in this SDG?  
 Y/N N/A Were target analytes detected in the field blanks?  
**Blank units:** mg/L Associated sample units: mg/L  
**Sampling date:** 8-2-10 Soil factor applied: NA  
**Field blank type:** (circle one) Field Blank / Rinsate / Other: Associated Samples: 1-9 (>RL)

Analyte	Blank ID	Action Level	No Qual's.	Sample Identification			
	FB080210V						
Cr	0.006						

**Blank units:** mg/L Associated sample units: mg/L  
**Sampling date:** 8-3-10 Soil factor applied: NA  
**Field blank type:** (circle one) Field Blank / Rinsate / Other: EB Associated Samples: 10-15 (>RL)

Analyte	Blank ID	Action Level	No Qual's.	Sample Identification			
	16						
Cr	0.0014						

CIRCLED RESULTS WERE NOT QUALIFIED. ALL RESULTS NOT CIRCLED WERE QUALIFIED BY THE FOLLOWING STATEMENT:  
 Samples with analyte concentrations within five times the associated field blank concentration are listed above, these sample results were qualified as not detected, "U".

LDC#: 2420304

**VALIDATION FINDINGS WORKSHEET**  
**Field Duplicates**

Page: 1 of 1  
Reviewer: MG  
2nd Reviewer: [Signature]

**METHOD:** Metals (EPA Method 6010B/6020/7000)

- Y N NA Were field duplicate pairs identified in this SDG?  
 Y N NA Were target analytes detected in the field duplicate pairs?

Analyte	Concentration (mg/L)		( $\leq 30$ )	(mg/L)	(mg/L)	Qualifications (Parent Only)
	10	15	RPD	Difference	Limits	
Chromium	0.11	0.098	12			

V:\FIELD DUPLICATES\FD\_inorganic\2420304.wpd

**Laboratory Data Consultants, Inc.  
Data Validation Report**

**Project/Site Name:** 2010 Annual Remedial Performance Sampling  
**Collection Date:** July 5, 2010  
**LDC Report Date:** December 19, 2010  
**Matrix:** Water  
**Parameters:** Wet Chemistry  
**Validation Level:** Stage 2A  
**Laboratory:** MWH Laboratories, Inc.  
**Sample Delivery Group (SDG):** 337662

**Sample Identification**

ART-1  
ART-2  
ART-3  
ART-4  
ART-6  
ART-7  
ART-8  
PC-99R2/R3  
PC-115R  
PC-116R  
SF-1  
PC-117  
PC-118  
PC-119  
PC-120  
PC-121  
PC-133  
ART-9

## Introduction

This data review covers 18 water samples listed on the cover sheet including dilutions and reanalysis as applicable. The analyses were per EPA Method 314.0 for Perchlorate, and EPA Method 160.1 and Standard Method 2540C for Total Dissolved Solids.

This review follows a modified outline of the USEPA Contract Laboratory Program National Functional Guidelines for Inorganic Data Review (October 2004) and the EPA Region 9 Superfund Data Evaluation/Validation Guidance, NDEP guidance (May 2006).

A qualification summary table is provided at the end of this report if data has been qualified. Flags are classified as P (protocol) or A (advisory) to indicate whether the flag is due to a laboratory deviation from a specified protocol or is of technical advisory nature.

Raw data were not reviewed for this SDG. The review was based on QC data.

The following are definitions of the data qualifiers:

- J+ Data are qualified as estimated, with a high bias likely to occur. False positives or false negatives are unlikely to have been reported.
- J- Data are qualified as estimated, with a low bias likely to occur. False positives or false negatives are unlikely to have been reported.
- J Data are qualified as estimated; it is not possible to assess the direction of the potential bias. False positives or false negatives are unlikely to have been reported.
- U Indicates the compound or analyte was analyzed for but not detected at or above the stated limit.
- R Data are qualified as rejected. There is a significant potential for the reporting of false negatives or false positives.
- UJ Indicates the compound or analyte was analyzed for but not detected. The sample detection limit is an estimated value.
- A Indicates the finding is based upon technical validation criteria.
- P Indicates the finding is related to a protocol/contractual deviation.
- None Indicates the data was not significantly impacted by the finding, therefore qualification was not required.

## **I. Technical Holding Times**

All technical holding time requirements were met.

The chain-of-custodies were reviewed for documentation of cooler temperatures. All cooler temperatures met validation criteria.

## **II. Calibration**

### **a. Initial Calibration**

Initial calibration data were not reviewed for Stage 2A.

### **b. Calibration Verification**

Calibration verification data were not reviewed for Stage 2A.

## **III. Blanks**

Method blanks were reviewed for each matrix as applicable. No contaminant concentrations were found in the preparation blanks.

No field blanks were identified in this SDG.

## **IV. Matrix Spike/Matrix Spike Duplicates**

The laboratory has indicated that there were no matrix spike (MS) analyses specified for the samples in this SDG, and therefore matrix spike analyses were not performed for this SDG.

## **V. Duplicates**

The laboratory has indicated that there were no duplicate (DUP) analyses specified for the samples in this SDG, and therefore duplicate analyses were not performed for this SDG.

## **VI. Laboratory Control Samples**

Laboratory control samples were reviewed for each matrix as applicable. Percent recoveries (%R) and relative percent differences (RPD) were within QC limits.

## **VII. Sample Result Verification**

Raw data were not reviewed for this SDG.

## **VIII. Overall Assessment of Data**

Data flags are summarized at the end of this report if data has been qualified.

## **IX. Field Duplicates**

No field duplicates were identified in this SDG.

**2010 Annual Remedial Performance Sampling  
Wet Chemistry - Data Qualification Summary - SDG 337662**

No Sample Data Qualified in this SDG

**2010 Annual Remedial Performance Sampling  
Wet Chemistry - Laboratory Blank Data Qualification Summary - SDG 337662**

No Sample Data Qualified in this SDG

**2010 Annual Remedial Performance Sampling  
Wet Chemistry - Field Blank Data Qualification Summary - SDG 337662**

No Sample Data Qualified in this SDG



Tronox Northgate Henderson

VALIDATION COMPLETENESS WORKSHEET

LDC #: 24203A6

SDG #: 337662

Laboratory: MWH Laboratories

Stage 2 ~~B~~ A

mg

Date: 12-1-10

Page: 1 of 1

Reviewer: MG

2nd Reviewer: [Signature]

METHOD: (Analyte) Perchlorate (EPA Method 314.0), TDS (EPA Method 160.1/SM2540C)

The samples listed below were reviewed for each of the following validation areas. Validation findings are noted in attached validation findings worksheets.

	Validation Area		Comments
I.	Technical holding times	A	Sampling dates: 7-5-10
Ia.	Initial calibration	N	
Iib.	Calibration verification	N	
III.	Blanks	A	
IV	Matrix Spike/Matrix Spike Duplicates	N	client specified
V	Duplicates	N	" "
VI.	Laboratory control samples	A	LCS/LCSD
VII.	Sample result verification	N	
VIII.	Overall assessment of data	A	
IX.	Field duplicates	N	
X	Field blanks	N	

Note: A = Acceptable  
N = Not provided/applicable  
SW = See worksheet

ND = No compounds detected  
R = Rinsate  
FB = Field blank

D = Duplicate  
TB = Trip blank  
EB = Equipment blank

Validated Samples:  
all water

1	ART-1	11	SF-1	21	PBW1	31	
2	ART-2	12	PC-117	22	PBW2	32	
3	ART-3	13	PC-118	23	PBW3	33	
4	ART-4	14	PC-119	24		34	
5	ART-6	15	PC-120	25		35	
6	ART-7	16	PC-121	26		36	
7	ART-8	17	PC-133	27		37	
8	PC-99R2/R3	18	ART-9	28		38	
9	PC-115R	19		29		39	
10	PC-116R	20		30		40	

Notes: \_\_\_\_\_

\_\_\_\_\_

\_\_\_\_\_

**VALIDATION FINDINGS WORKSHEET**  
**Sample Specific Analysis Reference**

All circled methods are applicable to each sample.

Sample ID	Matrix	Parameter
1→18	W	pH (TDS) Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> (ClO <sub>4</sub> )
		pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
		pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
		pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
		pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
		pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
		pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
		pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
		pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
		pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
		pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
		pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
		pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
		pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
		pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
		pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
		pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
		pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
		pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
		pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
		pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
		pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
		pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
		pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
		pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
		pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
		pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
		pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
		pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
		pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
		pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
		pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>

Comments: \_\_\_\_\_

## Laboratory Data Consultants, Inc. Data Validation Report

**Project/Site Name:** 2010 Annual Remedial Performance Sampling  
**Collection Date:** July 13 through July 15, 2010  
**LDC Report Date:** December 19, 2010  
**Matrix:** Water  
**Parameters:** Wet Chemistry  
**Validation Level:** Stage 2A  
**Laboratory:** MWH Laboratories, Inc.  
**Sample Delivery Group (SDG):** 338619

### Sample Identification

M-83	I-AD
M-87	MW-K4
PC-98R	ARP-1
PC-86	ARP-2A
PC-90	ARP-3A
PC-56	ARP-4A
PC-58	ARP-5A
PC-59	ARP-6B
PC-60	ARP-7
PC-62	PC-53
PC-68	PC-103
PC-122	MW-K5
PC-91	M-83DUP
PC-97	PC-68DUP
PC-18	
PC-55	
PC-101R	
PC-144	
ART-7B	
I-AC	

## Introduction

This data review covers 34 water samples listed on the cover sheet including dilutions and reanalysis as applicable. The analyses were per EPA Method 314.0 for Perchlorate, and EPA Method 160.1 and Standard Method 2540C for Total Dissolved Solids.

This review follows a modified outline of the USEPA Contract Laboratory Program National Functional Guidelines for Inorganic Data Review (October 2004) and the EPA Region 9 Superfund Data Evaluation/Validation Guidance, NDEP guidance (May 2006).

A qualification summary table is provided at the end of this report if data has been qualified. Flags are classified as P (protocol) or A (advisory) to indicate whether the flag is due to a laboratory deviation from a specified protocol or is of technical advisory nature.

Raw data were not reviewed for this SDG. The review was based on QC data.

The following are definitions of the data qualifiers:

J+ Data are qualified as estimated, with a high bias likely to occur. False positives or false negatives are unlikely to have been reported.

J- Data are qualified as estimated, with a low bias likely to occur. False positives or false negatives are unlikely to have been reported.

J Data are qualified as estimated; it is not possible to assess the direction of the potential bias. False positives or false negatives are unlikely to have been reported.

U Indicates the compound or analyte was analyzed for but not detected at or above the stated limit.

R Data are qualified as rejected. There is a significant potential for the reporting of false negatives or false positives.

UJ Indicates the compound or analyte was analyzed for but not detected. The sample detection limit is an estimated value.

A Indicates the finding is based upon technical validation criteria.

P Indicates the finding is related to a protocol/contractual deviation.

None Indicates the data was not significantly impacted by the finding, therefore qualification was not required.

## **I. Technical Holding Times**

All technical holding time requirements were met.

The chain-of-custodies were reviewed for documentation of cooler temperatures. All cooler temperatures met validation criteria.

## **II. Calibration**

### **a. Initial Calibration**

Initial calibration data were not reviewed for Stage 2A.

### **b. Calibration Verification**

Calibration verification data were not reviewed for Stage 2A.

## **III. Blanks**

Method blanks were reviewed for each matrix as applicable. No contaminant concentrations were found in the preparation blanks.

No field blanks were identified in this SDG.

## **IV. Matrix Spike/Matrix Spike Duplicates**

The laboratory has indicated that there were no matrix spike (MS) analyses specified for the samples in this SDG, and therefore matrix spike analyses were not performed for this SDG.

## **V. Duplicates**

Duplicate (DUP) sample analyses were reviewed for each matrix as applicable.

## **VI. Laboratory Control Samples**

Laboratory control samples were reviewed for each matrix as applicable. Percent recoveries (%R) and relative percent differences (RPD) were within QC limits.

## **VII. Sample Result Verification**

Raw data were not reviewed for this SDG.

## **VIII. Overall Assessment of Data**

Data flags are summarized at the end of this report if data has been qualified.

## **IX. Field Duplicates**

No field duplicates were identified in this SDG.

**2010 Annual Remedial Performance Sampling  
Wet Chemistry - Data Qualification Summary - SDG 338619**

No Sample Data Qualified in this SDG

**2010 Annual Remedial Performance Sampling  
Wet Chemistry - Laboratory Blank Data Qualification Summary - SDG 338619**

No Sample Data Qualified in this SDG

**2010 Annual Remedial Performance Sampling  
Wet Chemistry - Field Blank Data Qualification Summary - SDG 338619**

No Sample Data Qualified in this SDG

Tronox Northgate Henderson

VALIDATION COMPLETENESS WORKSHEET

LDC #: 24203B6

SDG #: 338619

Laboratory: MWH Laboratories

Stage 2B A

MM

Date: 7-1-10

Page: 1 of 1

Reviewer: MG

2nd Reviewer: [Signature]

METHOD: (Analyte) Perchlorate (EPA Method 314.0), TDS (EPA Method 160.1/SM2540C)

The samples listed below were reviewed for each of the following validation areas. Validation findings are noted in attached validation findings worksheets.

	Validation Area		Comments
I.	Technical holding times	A	Sampling dates: 7-13-10 through 7-15-10
IIa.	Initial calibration	N	
IIb.	Calibration verification	N	
III.	Blanks	A	
IV	Matrix Spike/Matrix Spike Duplicates	N	client specified
V	Duplicates	A	DUP
VI.	Laboratory control samples	A	LCS/LCSD
VII.	Sample result verification	N	
VIII.	Overall assessment of data	A	
IX.	Field duplicates	N	
X	Field blanks	N	

Note: A = Acceptable  
 N = Not provided/applicable  
 SW = See worksheet

ND = No compounds detected  
 R = Rinsate  
 FB = Field blank

D = Duplicate  
 TB = Trip blank  
 EB = Equipment blank

Validated Samples:  
 all water

1	M-83	11	PC-68	21	I-AD	31	PC-103
2	M-87	12	PC-122	22	MW-K4	32	MW-K5
3	PC-98R	13	PC-91	23	ARP-1	33	M-83DUP
4	PC-86	14	PC-97	24	ARP-2A	34	PC-68DUP
5	PC-90	15	PC-18	25	ARP-3A	35	
6	PC-56	16	PC-55	26	ARP-4A	36	
7	PC-58	17	PC-101R	27	ARP-5A	37	PBW1
8	PC-59	18	PC-144	28	ARP-6B	38	PBW2
9	PC-60	19	ART-7B	29	ARP-7	39	PBW3
10	PC-62	20	I-AC	30	PC-53	40	PBW4

Notes: \_\_\_\_\_



**VALIDATION FINDINGS WORKSHEET**  
**Sample Specific Analysis Reference**

All circled methods are applicable to each sample.

Sample ID	Matrix	Parameter
1 → 32	W	pH (TDS) CI F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> (ClO <sub>4</sub> )
ac 33,34	↓	pH (TDS) CI F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
		pH TDS CI F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
		pH TDS CI F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
		pH TDS CI F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
		pH TDS CI F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
		pH TDS CI F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
		pH TDS CI F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
		pH TDS CI F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
		pH TDS CI F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
		pH TDS CI F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
		pH TDS CI F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
		pH TDS CI F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
		pH TDS CI F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
		pH TDS CI F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
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Comments: \_\_\_\_\_

**Laboratory Data Consultants, Inc.  
Data Validation Report**

**Project/Site Name:** 2010 Annual Remedial Performance Sampling  
**Collection Date:** August 3 through August 4, 2010  
**LDC Report Date:** December 19, 2010  
**Matrix:** Water  
**Parameters:** Wet Chemistry  
**Validation Level:** Stage 2A  
**Laboratory:** MWH Laboratories, Inc.  
**Sample Delivery Group (SDG):** 340066

**Sample Identification**

M-99	M-88
M-25	I-V
M-92	VD080410
M-97	M-115DUP
M-14A	M-19MS
M-115	M-19MSD
M-17A	
M-34	
M-35	
M-19	
M-39	
I-K	
I-J	
I-AD	
I-Z	
I-I	
I-AC	
M-68	
M-74	
M-73	

## Introduction

This data review covers 26 water samples listed on the cover sheet including dilutions and reanalysis as applicable. The analyses were per EPA Method 314.0 for Perchlorate, and EPA Method 160.1 and Standard Method 2540C for Total Dissolved Solids.

This review follows a modified outline of the USEPA Contract Laboratory Program National Functional Guidelines for Inorganic Data Review (October 2004) and the EPA Region 9 Superfund Data Evaluation/Validation Guidance, NDEP guidance (May 2006).

A qualification summary table is provided at the end of this report if data has been qualified. Flags are classified as P (protocol) or A (advisory) to indicate whether the flag is due to a laboratory deviation from a specified protocol or is of technical advisory nature.

Raw data were not reviewed for this SDG. The review was based on QC data.

The following are definitions of the data qualifiers:

- J+ Data are qualified as estimated, with a high bias likely to occur. False positives or false negatives are unlikely to have been reported.
- J- Data are qualified as estimated, with a low bias likely to occur. False positives or false negatives are unlikely to have been reported.
- J Data are qualified as estimated; it is not possible to assess the direction of the potential bias. False positives or false negatives are unlikely to have been reported.
- U Indicates the compound or analyte was analyzed for but not detected at or above the stated limit.
- R Data are qualified as rejected. There is a significant potential for the reporting of false negatives or false positives.
- UJ Indicates the compound or analyte was analyzed for but not detected. The sample detection limit is an estimated value.
- A Indicates the finding is based upon technical validation criteria.
- P Indicates the finding is related to a protocol/contractual deviation.
- None Indicates the data was not significantly impacted by the finding, therefore qualification was not required.

## **I. Technical Holding Times**

All technical holding time requirements were met.

The chain-of-custodies were reviewed for documentation of cooler temperatures. All cooler temperatures met validation criteria.

## **II. Calibration**

### **a. Initial Calibration**

Initial calibration data were not reviewed for Stage 2A.

### **b. Calibration Verification**

Calibration verification data were not reviewed for Stage 2A.

## **III. Blanks**

Method blanks were reviewed for each matrix as applicable. No contaminant concentrations were found in the preparation blanks.

No field blanks were identified in this SDG.

## **IV. Matrix Spike/Matrix Spike Duplicates**

Matrix spike (MS) and matrix spike duplicate (MSD) samples were reviewed for each matrix as applicable. Percent recoveries (%R) and relative percent differences (RPD) were within QC limits.

## **V. Duplicates**

Duplicate (DUP) sample analyses were reviewed for each matrix as applicable. Results were within QC limits.

## **VI. Laboratory Control Samples**

Laboratory control samples were reviewed for each matrix as applicable. Percent recoveries (%R) and relative percent differences (RPD) were within QC limits.

## **VII. Sample Result Verification**

Raw data were not reviewed for this SDG.

## **VIII. Overall Assessment of Data**

Data flags are summarized at the end of this report if data has been qualified.

## IX. Field Duplicates

Samples M-19 and VD080410 were identified as field duplicates. No contaminant concentrations were detected in any of the samples with the following exceptions:

Analyte	Concentration		RPD (Limits)	Difference (Limits)	Flags	A or P
	M-121	ND-6				
Perchlorate	1600 ug/L	1700 ug/L	-	100 (≤400)	-	-
Total dissolved solids	3900 mg/L	3900 mg/L	0 (≤30)	-	-	-

**2010 Annual Remedial Performance Sampling  
Wet Chemistry - Data Qualification Summary - SDG 340066**

No Sample Data Qualified in this SDG

**2010 Annual Remedial Performance Sampling  
Wet Chemistry - Laboratory Blank Data Qualification Summary - SDG 340066**

No Sample Data Qualified in this SDG

**2010 Annual Remedial Performance Sampling  
Wet Chemistry - Field Blank Data Qualification Summary - SDG 340066**

No Sample Data Qualified in this SDG

LDC #: 24203C6  
 SDG #: 340066  
 Laboratory: MWH Laboratories

**Tronox Northgate Henderson**  
**VALIDATION COMPLETENESS WORKSHEET**  
 Stage 2 ~~B~~ **A**  
 gmA

Date: 8-1-10  
 Page: 1 of 1  
 Reviewer: MG  
 2nd Reviewer:           

**METHOD: (Analyte)** Perchlorate (EPA Method 314.0), TDS (EPA Method 160.1/SM2540C)

The samples listed below were reviewed for each of the following validation areas. Validation findings are noted in attached validation findings worksheets.

Validation Area			Comments
I.	Technical holding times	A	Sampling dates: 8-3-10 through 8-4-10
IIa.	Initial calibration	N	
IIb.	Calibration verification	N	
III.	Blanks	A	
IV.	Matrix Spike/Matrix Spike Duplicates	A	MS/MSD (SDG: 340161)
V.	Duplicates	A	DUP ( ↓ )
VI.	Laboratory control samples	A	LCS/LCSD
VII.	Sample result verification	N	
VIII.	Overall assessment of data	A	
IX.	Field duplicates	SW	D=10+23
X.	Field blanks	N	

Note: A = Acceptable  
 N = Not provided/applicable  
 SW = See worksheet  
 ND = No compounds detected  
 R = Rinsate  
 FB = Field blank  
 D = Duplicate  
 TB = Trip blank  
 EB = Equipment blank

Validated Samples: all water

1	M-99	11	M-39	21	M-88	31	
2	M-25	12	I-K	22	I-V	32	
3	M-92	13	I-J	23	VD080410	33	
4	M-97	14	I-AD	24	M-115DUP	34	
5	M-14A	15	I-Z	25	M-19MS	35	
6	M-115	16	I-I	26	M-19MSD	36	
7	M-17A	17	I-AC	27	PBW1	37	
8	M-34	18	M-68	28	PBW2	38	
9	M-35	19	M-74	29	PBW3	39	
10	M-19	20	M-73	30	PBW4	40	

Notes: \_\_\_\_\_  
 \_\_\_\_\_  
 \_\_\_\_\_

**VALIDATION FINDINGS WORKSHEET**  
**Sample Specific Analysis Reference**

All circled methods are applicable to each sample.

Sample ID	Matrix	Parameter
1 → 23	W	pH (TDS) CI F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> (ClO <sub>4</sub> )
QC ↓ 24	↓	pH (TDS) CI F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
↓ 25, 26	↓	pH TDS CI F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> (ClO <sub>4</sub> )
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Comments: \_\_\_\_\_  
 \_\_\_\_\_  
 \_\_\_\_\_



LDC#: 24203C6

SDG#: See Cover

### VALIDATION FINDINGS WORKSHEET

#### Field Duplicates

Page: 1 of 1

Reviewer: MG

2nd Reviewer: [Signature]

Inorganics, Method See Cover

N NA Were field duplicate pairs identified in this SDG?

N NA Were target analytes detected in the field duplicate pairs?

Analyte	Concentration (mg/L)		RPD ( $\leq 30$ )	Difference	Limits	Qualification (Parent only)
	10	23				
Perchlorate (ug/L)	1600	1700		100	( $\leq 400$ )	
TDS	3900	3900	0			

V:\FIELD DUPLICATES\FD\_inorganic\24203C6.wpd

**Laboratory Data Consultants, Inc.  
Data Validation Report**

**Project/Site Name:** 2010 Annual Remedial Performance Sampling  
**Collection Date:** August 4, 2010  
**LDC Report Date:** December 19, 2010  
**Matrix:** Water  
**Parameters:** Wet Chemistry  
**Validation Level:** Stage 4  
**Laboratory:** MWH Laboratories, Inc.  
**Sample Delivery Group (SDG):** 340161

**Sample Identification**

ART-1	PC-118DUP
ART-2	PC-119DUP
ART-3	
ART-4	
ART-6	
ART-7	
ART-8	
PC-99R2/R3	
PC-115R	
PC-116R	
PC-117	
PC-118	
PC-119	
PC-120	
PC-121	
PC-133	
ART-9	
ART-1MS	
ART-1MSD	
PC-117DUP	

## Introduction

This data review covers 22 water samples listed on the cover sheet including dilutions and reanalysis as applicable. The analyses were per EPA Method 314.0 for Perchlorate and EPA Method 160.1 and Standard Method 2540C for Total Dissolved Solids.

This review follows a modified outline of the USEPA Contract Laboratory Program National Functional Guidelines for Inorganic Data Review (October 2004) and the EPA Region 9 Superfund Data Evaluation/Validation Guidance, NDEP guidance (May 2006) as there are no current guidelines for the methods stated above.

A qualification summary table is provided at the end of this report if data has been qualified. Flags are classified as P (protocol) or A (advisory) to indicate whether the flag is due to a laboratory deviation from a specified protocol or is of technical advisory nature.

The following are definitions of the data qualifiers:

J+ Data are qualified as estimated, with a high bias likely to occur. False positives or false negatives are unlikely to have been reported.

J- Data are qualified as estimated, with a low bias likely to occur. False positives or false negatives are unlikely to have been reported.

J Data are qualified as estimated; it is not possible to assess the direction of the potential bias. False positives or false negatives are unlikely to have been reported.

U Indicates the compound or analyte was analyzed for but not detected at or above the stated limit.

R Data are qualified as rejected. There is a significant potential for the reporting of false negatives or false positives.

UJ Indicates the compound or analyte was analyzed for but not detected. The sample detection limit is an estimated value.

A Indicates the finding is based upon technical validation criteria.

P Indicates the finding is related to a protocol/contractual deviation.

None Indicates the data was not significantly impacted by the finding, therefore qualification was not required.

## **I. Technical Holding Times**

All technical holding time requirements were met.

The chain-of-custodies were reviewed for documentation of cooler temperatures. All cooler temperatures met validation criteria.

## **II. Calibration**

### **a. Initial Calibration**

All criteria for the initial calibration of each method were met.

### **b. Calibration Verification**

Calibration verification frequency and analysis criteria were met for each method when applicable.

## **III. Blanks**

Method blanks were reviewed for each matrix as applicable. No contaminant concentrations were found in the initial, continuing and preparation blanks.

No field blanks were identified in this SDG.

## **IV. Matrix Spike/Matrix Spike Duplicates**

Matrix spike (MS) and matrix spike duplicate (MSD) samples were reviewed for each matrix as applicable. Percent recoveries (%R) and relative percent differences (RPD) were within QC limits.

## **V. Duplicates**

Duplicate (DUP) sample analyses were reviewed for each matrix as applicable. Results were within QC limits.

## **VI. Laboratory Control Samples**

Laboratory control samples were reviewed for each matrix as applicable. Percent recoveries (%R) and relative percent differences (RPD) were within QC limits.

## **VII. Sample Result Verification**

All sample result verifications were acceptable.

## **VIII. Overall Assessment of Data**

Data flags are summarized at the end of this report if data has been qualified.

## **IX. Field Duplicates**

No field duplicates were identified in this SDG.

**2010 Annual Remedial Performance Sampling  
Wet Chemistry - Data Qualification Summary - SDG 340161**

No Sample Data Qualified in this SDG

**2010 Annual Remedial Performance Sampling  
Wet Chemistry - Laboratory Blank Data Qualification Summary - SDG 340161**

No Sample Data Qualified in this SDG

**2010 Annual Remedial Performance Sampling  
Wet Chemistry - Field Blank Data Qualification Summary - SDG 340161**

No Sample Data Qualified in this SDG

**Tronox Northgate Henderson**

**VALIDATION COMPLETENESS WORKSHEET**

LDC #: 24203D6

SDG #: 340161

Laboratory: MWH Laboratories

Stage 4

Date: 12-1-10

Page: 1 of 1

Reviewer: MG

2nd Reviewer: ✓

**METHOD: (Analyte)** Perchlorate (EPA Method 314.0), TDS (EPA Method 160.1/SM2540C)

The samples listed below were reviewed for each of the following validation areas. Validation findings are noted in attached validation findings worksheets.

	<b>Validation Area</b>		<b>Comments</b>
I.	Technical holding times	A	Sampling dates: <u>8-4-10</u>
IIa.	Initial calibration	A	
IIb.	Calibration verification	A	
III.	Blanks	A	
IV	Matrix Spike/Matrix Spike Duplicates	A	<u>MS/MSD</u>
V	Duplicates	A	<u>DUP</u>
VI.	Laboratory control samples	A	<u>LCS/LCSD</u>
VII.	Sample result verification	A	
VIII.	Overall assessment of data	A	
IX.	Field duplicates	N	
X	Field blanks	N	

Note: A = Acceptable  
N = Not provided/applicable  
SW = See worksheet

ND = No compounds detected  
R = Rinsate  
FB = Field blank

D = Duplicate  
TB = Trip blank  
EB = Equipment blank

Validated Samples:  
oil water

1	ART-1	11	PC-117	21	PC-118DUP	31	
2	ART-2	12	PC-118	22	PC-119DUP	32	
3	ART-3	13	PC-119	23		33	
4	ART-4	14	PC-120	24		34	
5	ART-6	15	PC-121	25		35	
6	ART-7	16	PC-133	26		36	
7	ART-8	17	ART-9	27		37	
8	PC-99R2/R3	18	ART-1MS	28		38	
9	PC-115R	19	ART-1MSD	29	<u>PBW1</u>	39	
10	PC-116R	20	PC-117DUP	30	<u>PBW2</u>	40	

Notes: \_\_\_\_\_  
\_\_\_\_\_  
\_\_\_\_\_

Method: Inorganics (EPA Method see cover)

Validation Area	Yes	No	NA	Findings/Comments
<b>I. Technical holding times</b>				
All technical holding times were met.	✓			
Cooler temperature criteria was met.	✓			
<b>II. Calibration</b>				
Were all instruments calibrated daily, each set-up time?	✓			
Were the proper number of standards used?	✓			
Were all initial calibration correlation coefficients $\geq 0.995$ ?	✓			
Were all initial and continuing calibration verification %Rs within the 90-110% QC limits?	✓			
Were titrant checks performed as required? (Level IV only)			✓	
Were balance checks performed as required? (Level IV only)	✓			
<b>III. Blanks</b>				
Was a method blank associated with every sample in this SDG?	✓			
Was there contamination in the method blanks? If yes, please see the Blanks validation completeness worksheet.		✓		
<b>IV. Matrix spike/Matrix spike duplicates and Duplicates</b>				
Were a matrix spike (MS) and duplicate (DUP) analyzed for each matrix in this SDG? If no, indicate which matrix does not have an associated MS/MSD or MS/DUP. Soil / Water.	✓			
Were the MS/MSD percent recoveries (%R) and the relative percent differences (RPD) within the 75-125 QC limits? If the sample concentration exceeded the spike concentration by a factor of 4 or more, no action was taken.	✓			
Were the MS/MSD or duplicate relative percent differences (RPD) $\leq 20\%$ for waters and $\leq 35\%$ for soil samples? A control limit of $\leq \text{CRDL}$ ( $\leq 2\text{X CRDL}$ for soil) was used for samples that were $\leq 5\text{X}$ the CRDL, including when only one of the duplicate sample values were $\leq 5\text{X}$ the CRDL.	✓			
<b>V. Laboratory control samples</b>				
Was an LCS analyzed for this SDG?	✓			
Was an LCS analyzed per extraction batch?	✓			
Were the LCS percent recoveries (%R) and relative percent difference (RPD) within the 80-120% (85-115% for Method 300.0) QC limits?	✓			
<b>VI. Regional Quality Assurance and Quality Control</b>				
Were performance evaluation (PE) samples performed?		✓		
Were the performance evaluation (PE) samples within the acceptance limits?			✓	



Validation Area	Yes	No	NA	Findings/Comments
<b>VII. Sample Result Verification</b>				
Were RLs adjusted to reflect all sample dilutions and dry weight factors applicable to level IV validation?	✓			
Were detection limits < RL?	✓			
<b>VIII. Overall assessment of data</b>				
Overall assessment of data was found to be acceptable.	✓			
<b>IX. Field duplicates</b>				
Field duplicate pairs were identified in this SDG.		✓		
Target analytes were detected in the field duplicates.			✓	
<b>X. Field blanks</b>				
Field blanks were identified in this SDG.		✓		
Target analytes were detected in the field blanks.			✓	

**VALIDATION FINDINGS WORKSHEET  
Sample Specific Analysis Reference**

All circled methods are applicable to each sample.

Sample ID	Matrix	Parameter
1 → 17	W	pH (TDS) Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> (ClO <sub>4</sub> )
QC 18 19	↓	pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> (ClO <sub>4</sub> )
↓ 20 → 22	↓	pH (TDS) Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
		pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
		pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
		pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
		pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
		pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
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		pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
		pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
		pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
		pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
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		pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
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		pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
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		pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
		pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
		pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
		pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
		pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
		pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
		pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
		pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
		pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
		pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
		pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>

Comments: \_\_\_\_\_

LDC #: 24203D6

VALIDATION FINDINGS WORKSHEET  
Initial and Continuing Calibration Calculation Verification

Page: 1 of 1  
Reviewer: MG  
2nd Reviewer: [Signature]

METHOD: Inorganics, Method See Cover

The correlation coefficient (r) for the calibration of C104 was recalculated. Calibration date: 8-31-10

An initial or continuing calibration verification percent recovery (%R) was recalculated for each type of analysis using the following formula:

%R = Found x 100 / True  
Where, Found = concentration of each analyte measured in the analysis of the ICV or CCV solution  
True = concentration of each analyte in the ICV or CCV source

Type of Analysis	Analyte	Standard ID	Found (units)	True (units)	Recalculated		Acceptable (Y/N)
					r or %R	Reported r or %R	
Initial calibration	C104	Blank	0.0 (mg/L)	0	r = 0.999541	r = 0.999547	Y
		Standard 1	2.0 ( )	0.005			
		Standard 2	4.0 ( )	0.011			
		Standard 3	10.0 ( )	0.027			
		Standard 4	25.0 ( )	0.071			
		Standard 5	50.0 ( )	0.146			
		Standard 6	75.0 ( )	0.229			
Standard 7	100.0 ( )	0.312					
Calibration verification	C104	Essee CCV	25.154 (mg/L)	25. (mg/L)	101	not reported	
Calibration verification	-	-	-	-	-	-	-
Calibration verification	-	-	-	-	-	-	-

Comments: Refer to Calibration Verification findings worksheet for list of qualifications and associated samples when reported results do not agree within 10.0% of the recalculated results.

**VALIDATION FINDINGS WORKSHEET**  
**Level IV Recalculation Worksheet**

**METHOD:** Inorganics, Method see cover

Percent recoveries (%R) for a laboratory control sample and a matrix spike sample were recalculated using the following formula:

$$\%R = \frac{\text{Found}}{\text{True}} \times 100$$
 Where, Found = concentration of each analyte measured in the analysis of the sample. For the matrix spike calculation, Found = SSR (spiked sample result) - SR (sample result).  
 True = concentration of each analyte in the source.

A sample and duplicate relative percent difference (RPD) was recalculated using the following formula:

$$RPD = \frac{|S-D|}{(S+D)/2} \times 100$$
 Where, S = Original sample concentration  
 D = Duplicate sample concentration

Sample ID	Type of Analysis	Element	Found / S (units)	True / D (units)	Recalculated		Reported		Acceptable (Y/N)
					%R / RPD	%R / RPD	%R / RPD	%R / RPD	
LCSI	Laboratory control sample	TDS	170 (mg/L)	175 (mg/L)	97	97	97	97	Y
18	Matrix spike sample	C104	(SSR-SR) 506.7 (μg/L)	500 (μg/L) 25 μg	101	101	101	101	↓
20	Duplicate sample	TDS	3308 (mg/L)	3296 (mg/L)	0.36	0.36	0.36	0.36	↓

Comments: Refer to appropriate worksheet for list of qualifications and associated samples when reported results do not agree within 10.0% of the recalculated results.

**VALIDATION FINDINGS WORKSHEET**  
Sample Calculation Verification

METHOD: Inorganics, Method See cover

Please see qualifications below for all questions answered "N". Not applicable questions are identified as "N/A".

- N N/A Have results been reported and calculated correctly?
- N N/A Are results within the calibrated range of the instruments?
- N N/A Are all detection limits below the CRQL?

Compound (analyte) results for # 1, ClO<sub>4</sub> reported with a positive detect were recalculated and verified using the following equation:

Concentration =  $y = mx + b$   
 where  
 $m = 0.0027$   
 $b = 0.0000$   
 $dil = 20x$

Recalculation:  
 $ClO_4 \text{ mg/L} = \left( \frac{0.253}{0.0027} \right) \times 20 = 1874 \text{ mg/L}$

#	Sample ID	Analyte	Reported Concentration ( )	Calculated Concentration ( )	Acceptable (Y/N)
1	1	ClO <sub>4</sub>	1700 (mg/L)	1900 (mg/L)	Y
		TDS	6400 (mg/L)	6400 (mg/L)	↓
2	11	ClO <sub>4</sub>	4100 (mg/L)	4400 (mg/L)	↓
		TDS	3300 (mg/L)	3300 (mg/L)	↓

Note: \_\_\_\_\_

**Laboratory Data Consultants, Inc.  
Data Validation Report**

**Project/Site Name:** 2010 Annual Remedial Performance Sampling

**Collection Date:** August 5, 2010

**LDC Report Date:** December 23, 2010

**Matrix:** Water

**Parameters:** Total Dissolved Solids

**Validation Level:** Stage 2A

**Laboratory:** MWH Laboratories, Inc.

**Sample Delivery Group (SDG):** 340226

**Sample Identification**

M-5A

## Introduction

This data review covers one water sample listed on the cover sheet including dilutions and reanalysis as applicable. The analyses were per EPA Method 160.1 and Standard Method 2540C for Total Dissolved Solids.

This review follows a modified outline of the USEPA Contract Laboratory Program National Functional Guidelines for Inorganic Data Review (October 2004) and the EPA Region 9 Superfund Data Evaluation/Validation Guidance, NDEP guidance (May 2006).

A qualification summary table is provided at the end of this report if data has been qualified. Flags are classified as P (protocol) or A (advisory) to indicate whether the flag is due to a laboratory deviation from a specified protocol or is of technical advisory nature.

Raw data were not reviewed for this SDG. The review was based on QC data.

The following are definitions of the data qualifiers:

J+ Data are qualified as estimated, with a high bias likely to occur. False positives or false negatives are unlikely to have been reported.

J- Data are qualified as estimated, with a low bias likely to occur. False positives or false negatives are unlikely to have been reported.

J Data are qualified as estimated; it is not possible to assess the direction of the potential bias. False positives or false negatives are unlikely to have been reported.

U Indicates the compound or analyte was analyzed for but not detected at or above the stated limit.

R Data are qualified as rejected. There is a significant potential for the reporting of false negatives or false positives.

UJ Indicates the compound or analyte was analyzed for but not detected. The sample detection limit is an estimated value.

A Indicates the finding is based upon technical validation criteria.

P Indicates the finding is related to a protocol/contractual deviation.

None Indicates the data was not significantly impacted by the finding, therefore qualification was not required.

## **I. Technical Holding Times**

All technical holding time requirements were met.

The chain-of-custodies were reviewed for documentation of cooler temperatures. All cooler temperatures met validation criteria.

## **II. Calibration**

### **a. Initial Calibration**

Initial calibration data were not reviewed for Stage 2A.

### **b. Calibration Verification**

Calibration verification data were not reviewed for Stage 2A.

## **III. Blanks**

Method blanks were reviewed for each matrix as applicable. No total dissolved solids were found in the preparation blanks.

No field blanks were identified in this SDG.

## **IV. Matrix Spike/Matrix Spike Duplicates**

The laboratory has indicated that there were no matrix spike (MS) analyses specified for the samples in this SDG, and therefore matrix spike analyses were not performed for this SDG.

## **V. Duplicates**

Duplicate (DUP) sample analyses were reviewed for each matrix as applicable. Results were within QC limits.

## **VI. Laboratory Control Samples**

Laboratory control samples were reviewed for each matrix as applicable. Percent recoveries (%R) and relative percent differences (RPD) were within QC limits.

## **VII. Sample Result Verification**

Raw data were not reviewed for this SDG.

## **VIII. Overall Assessment of Data**

Data flags are summarized at the end of this report if data has been qualified.



## **IX. Field Duplicates**

No field duplicates were identified in this SDG.

**2010 Annual Remedial Performance Sampling  
Wet Chemistry - Data Qualification Summary - SDG 340226**

No Sample Data Qualified in this SDG

**2010 Annual Remedial Performance Sampling  
Wet Chemistry - Laboratory Blank Data Qualification Summary - SDG 340226**

No Sample Data Qualified in this SDG

**2010 Annual Remedial Performance Sampling  
Wet Chemistry - Field Blank Data Qualification Summary - SDG 340226**

No Sample Data Qualified in this SDG

Tronox Northgate Henderson

VALIDATION COMPLETENESS WORKSHEET

LDC #: 24203E6

SDG #: 340226

Laboratory: MWH Laboratories

Stage 2 ~~B~~ A

gmf

Date: 12-1-10

Page: 1 of 1

Reviewer: MG

2nd Reviewer:

METHOD: (Analyte) TDS (EPA Method 160.1/SM2540C)

The samples listed below were reviewed for each of the following validation areas. Validation findings are noted in attached validation findings worksheets.

	Validation Area		Comments
I.	Technical holding times	A	Sampling dates: 8-5-10
IIa.	Initial calibration	N	
IIb.	Calibration verification	N	
III.	Blanks	A	
IV	Matrix Spike/Matrix Spike Duplicates	N	client specified
V	Duplicates	A	DUP (SDG: 340161)
VI.	Laboratory control samples	A	LCS/LCSD
VII.	Sample result verification	N	
VIII.	Overall assessment of data	A	
IX.	Field duplicates	N	
X	Field blanks	N	

Note: A = Acceptable  
 N = Not provided/applicable  
 SW = See worksheet  
 ND = No compounds detected  
 R = Rinsate  
 FB = Field blank  
 D = Duplicate  
 TB = Trip blank  
 EB = Equipment blank

Validated Samples: water

1	M-5A DUP 1	11		21		31	
2	PBW	12		22		32	
3		13		23		33	
4		14		24		34	
5		15		25		35	
6		16		26		36	
7		17		27		37	
8		18		28		38	
9		19		29		39	
10		20		30		40	

Notes: \_\_\_\_\_  
 \_\_\_\_\_  
 \_\_\_\_\_

**Laboratory Data Consultants, Inc.  
Data Validation Report**

**Project/Site Name:** 2010 Annual Remedial Performance Sampling  
**Collection Date:** August 5, 2010  
**LDC Report Date:** December 19, 2010  
**Matrix:** Water  
**Parameters:** Total Dissolved Solids  
**Validation Level:** Stage 2A  
**Laboratory:** MWH Laboratories, Inc.  
**Sample Delivery Group (SDG):** 340229

**Sample Identification**

M7B

## Introduction

This data review covers one water sample listed on the cover sheet including dilutions and reanalysis as applicable. The analyses were per EPA Method 160.1 and Standard Method 2540C for Total Dissolved Solids.

This review follows a modified outline of the USEPA Contract Laboratory Program National Functional Guidelines for Inorganic Data Review (October 2004) and the EPA Region 9 Superfund Data Evaluation/Validation Guidance, NDEP guidance (May 2006).

A qualification summary table is provided at the end of this report if data has been qualified. Flags are classified as P (protocol) or A (advisory) to indicate whether the flag is due to a laboratory deviation from a specified protocol or is of technical advisory nature.

Raw data were not reviewed for this SDG. The review was based on QC data.

The following are definitions of the data qualifiers:

- J+ Data are qualified as estimated, with a high bias likely to occur. False positives or false negatives are unlikely to have been reported.
- J- Data are qualified as estimated, with a low bias likely to occur. False positives or false negatives are unlikely to have been reported.
- J Data are qualified as estimated; it is not possible to assess the direction of the potential bias. False positives or false negatives are unlikely to have been reported.
- U Indicates the compound or analyte was analyzed for but not detected at or above the stated limit.
- R Data are qualified as rejected. There is a significant potential for the reporting of false negatives or false positives.
- UU Indicates the compound or analyte was analyzed for but not detected. The sample detection limit is an estimated value.
- A Indicates the finding is based upon technical validation criteria.
- P Indicates the finding is related to a protocol/contractual deviation.
- None Indicates the data was not significantly impacted by the finding, therefore qualification was not required.

## **I. Technical Holding Times**

All technical holding time requirements were met.

The chain-of-custodies were reviewed for documentation of cooler temperatures. All cooler temperatures met validation criteria.

## **II. Calibration**

### **a. Initial Calibration**

Initial calibration data were not reviewed for Stage 2A.

### **b. Calibration Verification**

Calibration verification data were not reviewed for Stage 2A.

## **III. Blanks**

Method blanks were reviewed for each matrix as applicable. No total dissolved solids were found in the preparation blanks.

No field blanks were identified in this SDG.

## **IV. Matrix Spike/Matrix Spike Duplicates**

The laboratory has indicated that there were no matrix spike (MS) analyses specified for the samples in this SDG, and therefore matrix spike analyses were not performed for this SDG.

## **V. Duplicates**

Duplicate (DUP) sample analyses were reviewed for each matrix as applicable. Results were within QC limits.

## **VI. Laboratory Control Samples**

Laboratory control samples were reviewed for each matrix as applicable. Percent recoveries (%R) and relative percent differences (RPD) were within QC limits.

## **VII. Sample Result Verification**

Raw data were not reviewed for this SDG.

## **VIII. Overall Assessment of Data**

Data flags are summarized at the end of this report if data has been qualified.

## **IX. Field Duplicates**

No field duplicates were identified in this SDG.

**2010 Annual Remedial Performance Sampling  
Wet Chemistry - Data Qualification Summary - SDG 340229**

No Sample Data Qualified in this SDG

**2010 Annual Remedial Performance Sampling  
Wet Chemistry - Laboratory Blank Data Qualification Summary - SDG 340229**

No Sample Data Qualified in this SDG

**2010 Annual Remedial Performance Sampling  
Wet Chemistry - Field Blank Data Qualification Summary - SDG 340229**

No Sample Data Qualified in this SDG



Tronox Northgate Henderson

LDC #: 24203F6

VALIDATION COMPLETENESS WORKSHEET

Date: 12-1-10

SDG #: 340229

Stage 2B A

Page: 1 of 1

Laboratory: MWH Laboratories

9M4

Reviewer: MG

2nd Reviewer: ✓

METHOD: (Analyte) TDS (EPA Method 160.1/SM2540C)

The samples listed below were reviewed for each of the following validation areas. Validation findings are noted in attached validation findings worksheets.

	Validation Area		Comments
I.	Technical holding times	A	Sampling dates: 8-5-10
IIa.	Initial calibration	N	
IIb.	Calibration verification	N	
III.	Blanks	A	
IV	Matrix Spike/Matrix Spike Duplicates	N	client specified
V	Duplicates	A	DUP (SDG: 340161)
VI.	Laboratory control samples	A	LCS/LCSD
VII.	Sample result verification	N	
VIII.	Overall assessment of data	A	
IX.	Field duplicates	N	
X	Field blanks	N	

Note: A = Acceptable  
N = Not provided/applicable  
SW = See worksheet

ND = No compounds detected  
R = Rinsate  
FB = Field blank

D = Duplicate  
TB = Trip blank  
EB = Equipment blank

Validated Samples:  
Water

1	M7B DUP1	11		21		31	
2	PBW	12		22		32	
3		13		23		33	
4		14		24		34	
5		15		25		35	
6		16		26		36	
7		17		27		37	
8		18		28		38	
9		19		29		39	
10		20		30		40	

Notes: \_\_\_\_\_  
\_\_\_\_\_  
\_\_\_\_\_

**Laboratory Data Consultants, Inc.  
Data Validation Report**

**Project/Site Name:** 2010 Annual Remedial Performance Sampling  
**Collection Date:** August 6, 2010  
**LDC Report Date:** December 19, 2010  
**Matrix:** Water  
**Parameters:** Total Dissolved Solids  
**Validation Level:** Stage 2A  
**Laboratory:** MWH Laboratories, Inc.  
**Sample Delivery Group (SDG):** 340275

**Sample Identification**

H-28A

## Introduction

This data review covers one water sample listed on the cover sheet including dilutions and reanalysis as applicable. The analyses were per EPA Method 160.1 and Standard Method 2540C for Total Dissolved Solids.

This review follows a modified outline of the USEPA Contract Laboratory Program National Functional Guidelines for Inorganic Data Review (October 2004) and the EPA Region 9 Superfund Data Evaluation/Validation Guidance, NDEP guidance (May 2006).

A qualification summary table is provided at the end of this report if data has been qualified. Flags are classified as P (protocol) or A (advisory) to indicate whether the flag is due to a laboratory deviation from a specified protocol or is of technical advisory nature.

Raw data were not reviewed for this SDG. The review was based on QC data.

The following are definitions of the data qualifiers:

J+ Data are qualified as estimated, with a high bias likely to occur. False positives or false negatives are unlikely to have been reported.

J- Data are qualified as estimated, with a low bias likely to occur. False positives or false negatives are unlikely to have been reported.

J Data are qualified as estimated; it is not possible to assess the direction of the potential bias. False positives or false negatives are unlikely to have been reported.

U Indicates the compound or analyte was analyzed for but not detected at or above the stated limit.

R Data are qualified as rejected. There is a significant potential for the reporting of false negatives or false positives.

UJ Indicates the compound or analyte was analyzed for but not detected. The sample detection limit is an estimated value.

A Indicates the finding is based upon technical validation criteria.

P Indicates the finding is related to a protocol/contractual deviation.

None Indicates the data was not significantly impacted by the finding, therefore qualification was not required.

## **I. Technical Holding Times**

All technical holding time requirements were met.

The chain-of-custodies were reviewed for documentation of cooler temperatures. All cooler temperatures met validation criteria.

## **II. Calibration**

### **a. Initial Calibration**

Initial calibration data were not reviewed for Stage 2A.

### **b. Calibration Verification**

Calibration verification data were not reviewed for Stage 2A.

## **III. Blanks**

Method blanks were reviewed for each matrix as applicable. No total dissolved solids were found in the preparation blanks.

No field blanks were identified in this SDG.

## **IV. Matrix Spike/Matrix Spike Duplicates**

The laboratory has indicated that there were no matrix spike (MS) analyses specified for the samples in this SDG, and therefore matrix spike analyses were not performed for this SDG.

## **V. Duplicates**

The laboratory has indicated that there were no duplicate (DUP) analyses specified for the samples in this SDG, and therefore duplicate analyses were not performed for this SDG.

## **VI. Laboratory Control Samples**

Laboratory control samples were reviewed for each matrix as applicable. Percent recoveries (%R) and relative percent differences (RPD) were within QC limits.

## **VII. Sample Result Verification**

Raw data were not reviewed for this SDG.

## **VIII. Overall Assessment of Data**

Data flags are summarized at the end of this report if data has been qualified.

## **IX. Field Duplicates**

No field duplicates were identified in this SDG.

**2010 Annual Remedial Performance Sampling  
Wet Chemistry - Data Qualification Summary - SDG 340275**

No Sample Data Qualified in this SDG

**2010 Annual Remedial Performance Sampling  
Wet Chemistry - Laboratory Blank Data Qualification Summary - SDG 340275**

No Sample Data Qualified in this SDG

**2010 Annual Remedial Performance Sampling  
Wet Chemistry - Field Blank Data Qualification Summary - SDG 340275**

No Sample Data Qualified in this SDG

LDC #: 24203G6  
 SDG #: 340275  
 Laboratory: MWH Laboratories

**Tronox Northgate Henderson**  
**VALIDATION COMPLETENESS WORKSHEET**

Stage 2B A

*gmA*

Date: 12-1-10  
 Page: 1 of 1  
 Reviewer: MG  
 2nd Reviewer: [Signature]

**METHOD: (Analyte)** TDS (EPA Method 160.1/SM2540C)

The samples listed below were reviewed for each of the following validation areas. Validation findings are noted in attached validation findings worksheets.

	Validation Area		Comments
I.	Technical holding times	A	Sampling dates: <u>8-6-10</u>
IIa.	Initial calibration	N	
IIb.	Calibration verification	N	
III.	Blanks	A	
IV	Matrix Spike/Matrix Spike Duplicates	N	<u>client specified</u>
V	Duplicates	N	<u>" "</u>
VI.	Laboratory control samples	A	<u>LCS/LCSD</u>
VII.	Sample result verification	N	
VIII.	Overall assessment of data	A	
IX.	Field duplicates	N	
X	Field blanks	N	

Note: A = Acceptable      ND = No compounds detected      D = Duplicate  
 N = Not provided/applicable      R = Rinsate      TB = Trip blank  
 SW = See worksheet      FB = Field blank      EB = Equipment blank

Validated Samples:  
water

1	<u>H-28A DUPT</u>	11		21		31	
2	<u>PBW</u>	12		22		32	
3		13		23		33	
4		14		24		34	
5		15		25		35	
6		16		26		36	
7		17		27		37	
8		18		28		38	
9		19		29		39	
10		20		30		40	

Notes: \_\_\_\_\_  
 \_\_\_\_\_  
 \_\_\_\_\_

**Laboratory Data Consultants, Inc.  
Data Validation Report**

**Project/Site Name:** 2010 Annual Remedial Performance Sampling

**Collection Date:** August 6, 2010

**LDC Report Date:** December 19, 2010

**Matrix:** Water

**Parameters:** Total Dissolved Solids

**Validation Level:** Stage 2A

**Laboratory:** MWH Laboratories, Inc.

**Sample Delivery Group (SDG):** 340276

**Sample Identification**

M-6A



## Introduction

This data review covers one water sample listed on the cover sheet including dilutions and reanalysis as applicable. The analyses were per EPA Method 160.1 and Standard Method 2540C for Total Dissolved Solids.

This review follows a modified outline of the USEPA Contract Laboratory Program National Functional Guidelines for Inorganic Data Review (October 2004) and the EPA Region 9 Superfund Data Evaluation/Validation Guidance, NDEP guidance (May 2006).

A qualification summary table is provided at the end of this report if data has been qualified. Flags are classified as P (protocol) or A (advisory) to indicate whether the flag is due to a laboratory deviation from a specified protocol or is of technical advisory nature.

Raw data were not reviewed for this SDG. The review was based on QC data.

The following are definitions of the data qualifiers:

J+ Data are qualified as estimated, with a high bias likely to occur. False positives or false negatives are unlikely to have been reported.

J- Data are qualified as estimated, with a low bias likely to occur. False positives or false negatives are unlikely to have been reported.

J Data are qualified as estimated; it is not possible to assess the direction of the potential bias. False positives or false negatives are unlikely to have been reported.

U Indicates the compound or analyte was analyzed for but not detected at or above the stated limit.

R Data are qualified as rejected. There is a significant potential for the reporting of false negatives or false positives.

UJ Indicates the compound or analyte was analyzed for but not detected. The sample detection limit is an estimated value.

A Indicates the finding is based upon technical validation criteria.

P Indicates the finding is related to a protocol/contractual deviation.

None Indicates the data was not significantly impacted by the finding, therefore qualification was not required.

## **I. Technical Holding Times**

All technical holding time requirements were met.

The chain-of-custodies were reviewed for documentation of cooler temperatures. All cooler temperatures met validation criteria.

## **II. Calibration**

### **a. Initial Calibration**

Initial calibration data were not reviewed for Stage 2A.

### **b. Calibration Verification**

Calibration verification data were not reviewed for Stage 2A.

## **III. Blanks**

Method blanks were reviewed for each matrix as applicable. No total dissolved solids were found in the preparation blanks.

No field blanks were identified in this SDG.

## **IV. Matrix Spike/Matrix Spike Duplicates**

The laboratory has indicated that there were no matrix spike (MS) analyses specified for the samples in this SDG, and therefore matrix spike analyses were not performed for this SDG.

## **V. Duplicates**

The laboratory has indicated that there were no duplicate (DUP) analyses specified for the samples in this SDG, and therefore duplicate analyses were not performed for this SDG.

## **VI. Laboratory Control Samples**

Laboratory control samples were reviewed for each matrix as applicable. Percent recoveries (%R) and relative percent differences (RPD) were within QC limits.

## **VII. Sample Result Verification**

Raw data were not reviewed for this SDG.

## **VIII. Overall Assessment of Data**

Data flags are summarized at the end of this report if data has been qualified.

## **IX. Field Duplicates**

No field duplicates were identified in this SDG.

**2010 Annual Remedial Performance Sampling  
Wet Chemistry - Data Qualification Summary - SDG 340276**

No Sample Data Qualified in this SDG

**2010 Annual Remedial Performance Sampling  
Wet Chemistry - Laboratory Blank Data Qualification Summary - SDG 340276**

No Sample Data Qualified in this SDG

**2010 Annual Remedial Performance Sampling  
Wet Chemistry - Field Blank Data Qualification Summary - SDG 340276**

No Sample Data Qualified in this SDG

Tronox Northgate Henderson

VALIDATION COMPLETENESS WORKSHEET

LDC #: 24203H6  
 SDG #: 340276  
 Laboratory: MWH Laboratories

Stage 2B A

9M

Date: 10-1-10  
 Page: 1 of 1  
 Reviewer: MG  
 2nd Reviewer: V

METHOD: (Analyte) TDS (EPA Method 160.1/SM2540C)

The samples listed below were reviewed for each of the following validation areas. Validation findings are noted in attached validation findings worksheets.

	Validation Area		Comments
I.	Technical holding times	A	Sampling dates: <u>8-6-10</u>
IIa.	Initial calibration	N	
IIb.	Calibration verification	N	
III.	Blanks	A	
IV	Matrix Spike/Matrix Spike Duplicates	N	<u>client specified</u>
V	Duplicates	N	<u>" "</u>
VI.	Laboratory control samples	A	<u>LCS/LCSD</u>
VII.	Sample result verification	N	
VIII.	Overall assessment of data	A	
IX.	Field duplicates	N	
X	Field blanks	N	

Note: A = Acceptable  
 N = Not provided/applicable  
 SW = See worksheet  
 ND = No compounds detected  
 R = Rinsate  
 FB = Field blank  
 D = Duplicate  
 TB = Trip blank  
 EB = Equipment blank

Validated Samples:  
water

1	<u>M-6A DUP 1</u>	11		21		31	
2	<u>PBW</u>	12		22		32	
3		13		23		33	
4		14		24		34	
5		15		25		35	
6		16		26		36	
7		17		27		37	
8		18		28		38	
9		19		29		39	
10		20		30		40	

Notes: \_\_\_\_\_  
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**Laboratory Data Consultants, Inc.  
Data Validation Report**

**Project/Site Name:** 2010 Annual Remedial Performance Sampling  
**Collection Date:** August 6, 2010  
**LDC Report Date:** December 19, 2010  
**Matrix:** Water  
**Parameters:** Total Dissolved Solids  
**Validation Level:** Stage 2A  
**Laboratory:** MWH Laboratories, Inc.  
**Sample Delivery Group (SDG):** 340278

**Sample Identification**

M-10

## Introduction

This data review covers one water sample listed on the cover sheet including dilutions and reanalysis as applicable. The analyses were per EPA Method 160.1 and Standard Method 2540C for Total Dissolved Solids.

This review follows a modified outline of the USEPA Contract Laboratory Program National Functional Guidelines for Inorganic Data Review (October 2004) and the EPA Region 9 Superfund Data Evaluation/Validation Guidance, NDEP guidance (May 2006).

A qualification summary table is provided at the end of this report if data has been qualified. Flags are classified as P (protocol) or A (advisory) to indicate whether the flag is due to a laboratory deviation from a specified protocol or is of technical advisory nature.

Raw data were not reviewed for this SDG. The review was based on QC data.

The following are definitions of the data qualifiers:

- J+ Data are qualified as estimated, with a high bias likely to occur. False positives or false negatives are unlikely to have been reported.
- J- Data are qualified as estimated, with a low bias likely to occur. False positives or false negatives are unlikely to have been reported.
- J Data are qualified as estimated; it is not possible to assess the direction of the potential bias. False positives or false negatives are unlikely to have been reported.
- U Indicates the compound or analyte was analyzed for but not detected at or above the stated limit.
- R Data are qualified as rejected. There is a significant potential for the reporting of false negatives or false positives.
- UJ Indicates the compound or analyte was analyzed for but not detected. The sample detection limit is an estimated value.
- A Indicates the finding is based upon technical validation criteria.
- P Indicates the finding is related to a protocol/contractual deviation.
- None Indicates the data was not significantly impacted by the finding, therefore qualification was not required.

## **I. Technical Holding Times**

All technical holding time requirements were met.

The chain-of-custodies were reviewed for documentation of cooler temperatures. All cooler temperatures met validation criteria.

## **II. Calibration**

### **a. Initial Calibration**

Initial calibration data were not reviewed for Stage 2A.

### **b. Calibration Verification**

Calibration verification data were not reviewed for Stage 2A.

## **III. Blanks**

Method blanks were reviewed for each matrix as applicable. No total dissolved solids were found in the preparation blanks.

No field blanks were identified in this SDG.

## **IV. Matrix Spike/Matrix Spike Duplicates**

The laboratory has indicated that there were no matrix spike (MS) analyses specified for the samples in this SDG, and therefore matrix spike analyses were not performed for this SDG.

## **V. Duplicates**

The laboratory has indicated that there were no duplicate (DUP) analyses specified for the samples in this SDG, and therefore duplicate analyses were not performed for this SDG.

## **VI. Laboratory Control Samples**

Laboratory control samples were reviewed for each matrix as applicable. Percent recoveries (%R) and relative percent differences (RPD) were within QC limits.

## **VII. Sample Result Verification**

Raw data were not reviewed for this SDG.

## **VIII. Overall Assessment of Data**

Data flags are summarized at the end of this report if data has been qualified.



## **IX. Field Duplicates**

No field duplicates were identified in this SDG.

**2010 Annual Remedial Performance Sampling  
Wet Chemistry - Data Qualification Summary - SDG 340278**

No Sample Data Qualified in this SDG

**2010 Annual Remedial Performance Sampling  
Wet Chemistry - Laboratory Blank Data Qualification Summary - SDG 340278**

No Sample Data Qualified in this SDG

**2010 Annual Remedial Performance Sampling  
Wet Chemistry - Field Blank Data Qualification Summary - SDG 340278**

No Sample Data Qualified in this SDG

LDC #: 2420316  
 SDG #: 340278  
 Laboratory: MWH Laboratories

Tronox Northgate Henderson  
**VALIDATION COMPLETENESS WORKSHEET**

Stage 2 ~~B~~ A  
 mH

Date: 12-2-10  
 Page: 1 of 1  
 Reviewer: MG  
 2nd Reviewer:           

**METHOD: (Analyte)** TDS (EPA Method 160.1/SM2540C)

The samples listed below were reviewed for each of the following validation areas. Validation findings are noted in attached validation findings worksheets.

	Validation Area		Comments
I.	Technical holding times	A	Sampling dates: 8-6-10
IIa.	Initial calibration	<del>A</del> N	mH
IIb.	Calibration verification	N	
III.	Blanks	A	
IV	Matrix Spike/Matrix Spike Duplicates	N	client specified
V	Duplicates	N	" "
VI.	Laboratory control samples	A	LCS/LCSD
VII.	Sample result verification	N	
VIII.	Overall assessment of data	A	
IX.	Field duplicates	N	
X	Field blanks	N	

Note: A = Acceptable                      ND = No compounds detected                      D = Duplicate  
 N = Not provided/applicable                      R = Rinsate                      TB = Trip blank  
 SW = See worksheet                      FB = Field blank                      EB = Equipment blank

Validated Samples:  
 Water

1	M-10	11		21		31	
2	PBW	12		22		32	
3		13		23		33	
4		14		24		34	
5		15		25		35	
6		16		26		36	
7		17		27		37	
8		18		28		38	
9		19		29		39	
10		20		30		40	

Notes: \_\_\_\_\_  
 \_\_\_\_\_  
 \_\_\_\_\_

**Laboratory Data Consultants, Inc.  
Data Validation Report**

**Project/Site Name:** 2010 Annual Remedial Performance Sampling  
**Collection Date:** August 9 through August 12, 2010  
**LDC Report Date:** December 27, 2010  
**Matrix:** Water  
**Parameters:** Wet Chemistry  
**Validation Level:** Stage 2A  
**Laboratory:** MWH Laboratories, Inc.  
**Sample Delivery Group (SDG):** 340887

**Sample Identification**

M-83	MW-K5
M-87	PC-91
PC-86	PC-97
PC-90	PC-18
PC-56	PC-55
PC-58	PC-101R
PC-59	ART-7B
PC-60	PC-92
PC-62	PC-94
PC-68	M-83DUP
PC-122	PC-86DUP
MW-K4	MW-K5MS
ARP-1	MW-K5MSD
ARP-2A	
ARP-3A	
ARP-4A	
ARP-5A	
ARP-6B	
ARP-7	
PC-53	

## Introduction

This data review covers 33 water samples listed on the cover sheet including dilutions and reanalysis as applicable. The analyses were per EPA Method 314.0 for Perchlorate, and EPA Method 160.1 and Standard Method 2540C for Total Dissolved Solids.

This review follows a modified outline of the USEPA Contract Laboratory Program National Functional Guidelines for Inorganic Data Review (October 2004) and the EPA Region 9 Superfund Data Evaluation/Validation Guidance, NDEP guidance (May 2006).

A qualification summary table is provided at the end of this report if data has been qualified. Flags are classified as P (protocol) or A (advisory) to indicate whether the flag is due to a laboratory deviation from a specified protocol or is of technical advisory nature.

Raw data were not reviewed for this SDG. The review was based on QC data.

The following are definitions of the data qualifiers:

- J+ Data are qualified as estimated, with a high bias likely to occur. False positives or false negatives are unlikely to have been reported.
- J- Data are qualified as estimated, with a low bias likely to occur. False positives or false negatives are unlikely to have been reported.
- J Data are qualified as estimated; it is not possible to assess the direction of the potential bias. False positives or false negatives are unlikely to have been reported.
- U Indicates the compound or analyte was analyzed for but not detected at or above the stated limit.
- R Data are qualified as rejected. There is a significant potential for the reporting of false negatives or false positives.
- UJ Indicates the compound or analyte was analyzed for but not detected. The sample detection limit is an estimated value.
- A Indicates the finding is based upon technical validation criteria.
- P Indicates the finding is related to a protocol/contractual deviation.
- None Indicates the data was not significantly impacted by the finding, therefore qualification was not required.

## I. Technical Holding Times

All technical holding time requirements were met with the following exceptions:

Sample	Analyte	Total Time From Sample Collection Until Analysis	Required Holding Time From Sample Collection Until Analysis	Flag	A or P
PC-101R	Perchlorate	29 days	28 days	J- (all detects) UJ (all nondetects)	A

The chain-of-custodies were reviewed for documentation of cooler temperatures. All cooler temperatures met validation criteria.

## II. Calibration

### a. Initial Calibration

Initial calibration data were not reviewed for Stage 2A.

### b. Calibration Verification

Calibration verification data were not reviewed for Stage 2A.

## III. Blanks

Method blanks were reviewed for each matrix as applicable. No contaminant concentrations were found in the preparation blanks.

No field blanks were identified in this SDG.

## IV. Matrix Spike/Matrix Spike Duplicates

Matrix spike (MS) and matrix spike duplicate (MSD) samples were reviewed for each matrix as applicable. Percent recoveries (%R) and relative percent differences (RPD) were within QC limits.

## V. Duplicates

Duplicate (DUP) sample analyses were reviewed for each matrix as applicable. Results were within QC limits.

## VI. Laboratory Control Samples

Laboratory control samples were reviewed for each matrix as applicable. Percent recoveries (%R) and relative percent differences (RPD) were within QC limits with the following exceptions:

LCS ID (Associated Samples)	Analyte	LCS %R (Limits)	LCSD %R (Limits)	RPD (Limits)	Flag	A or P
LCS/LCSD (PC-90 PC-122 MW-K4 ARP-1 ARP-2A ARP-3A ARP-4A ARP-5A ARP-6B ARP-7 PC-53 MW-K5 PC-91 PC-97 PC-18 PC-92 PC-94)	Perchlorate	79 (85-115)	-	26 (≤20)	J (all detects) UJ (all non-detects)	P

### VII. Sample Result Verification

Raw data were not reviewed for this SDG.

### VIII. Overall Assessment of Data

Data flags are summarized at the end of this report if data has been qualified.

### IX. Field Duplicates

No field duplicates were identified in this SDG.

**2010 Annual Remedial Performance Sampling  
Wet Chemistry - Data Qualification Summary - SDG 340887**

SDG	Sample	Analyte	Flag	A or P	Reason
340887	PC-101R	Perchlorate	J- (all detects) UJ (all nondetects)	A	Technical holding times
340887	PC-90 PC-122 MW-K4 ARP-1 ARP-2A ARP-3A ARP-4A ARP-5A ARP-6B ARP-7 PC-53 MW-K5 PC-91 PC-97 PC-18 PC-92 PC-94	Perchlorate	J (all detects) UJ (all non-detects)	P	Laboratory control samples (%R)(RPD)

**2010 Annual Remedial Performance Sampling  
Wet Chemistry - Laboratory Blank Data Qualification Summary - SDG 340887**

No Sample Data Qualified in this SDG

**2010 Annual Remedial Performance Sampling  
Wet Chemistry - Field Blank Data Qualification Summary - SDG 340887**

No Sample Data Qualified in this SDG



**Tronox Northgate Henderson**

**VALIDATION COMPLETENESS WORKSHEET**

LDC #: 24203J6  
 SDG #: 340887  
 Laboratory: MWH Laboratories

Stage ~~2~~ **A**  
*gmf*

Date: 12-2-10  
 Page: 1 of 1  
 Reviewer: MCT  
 2nd Reviewer: [Signature]

**METHOD: (Analyte)** Perchlorate (EPA Method 314.0), TDS (EPA Method 160.1/SM2540C)

The samples listed below were reviewed for each of the following validation areas. Validation findings are noted in attached validation findings worksheets.

	<b>Validation Area</b>		<b>Comments</b>
I.	Technical holding times	SW	Sampling dates: <u>8-9-10 through 8-12-10</u>
IIa.	Initial calibration	N	
IIb.	Calibration verification	N	
III.	Blanks	A	
IV	Matrix Spike/Matrix Spike Duplicates	A	<u>MS/MSD</u>
V	Duplicates	A	<u>DUP</u>
VI.	Laboratory control samples	SW	<u>LCS/LCSD</u>
VII.	Sample result verification	N	
VIII.	Overall assessment of data	A	
IX.	Field duplicates	N	
X	Field blanks	N	

Note: A = Acceptable      ND = No compounds detected      D = Duplicate  
 N = Not provided/applicable      R = Rinsate      TB = Trip blank  
 SW = See worksheet      FB = Field blank      EB = Equipment blank

Validated Samples: all water

1	M-83	11	PC-122	21	MW-K5	31	PC-86DUP
2	M-87	12	MW-K4	22	PC-91	32	MW-K5MS
3	PC-86	13	ARP-1	23	PC-97	33	MW-K5MSD
4	PC-90	14	ARP-2A	24	PC-18	34	
5	PC-56	15	ARP-3A	25	PC-55	35	
6	PC-58	16	ARP-4A	26	PC-101R	36	
7	PC-59	17	ARP-5A	27	ART-7B	37	
8	PC-60	18	ARP-6B	28	PC-92	38	<u>PBW1</u>
9	PC-62	19	ARP-7	29	PC-94	39	<u>PBW2</u>
10	PC-68	20	PC-53	30	M-83DUP	40	<u>PBW3</u>

Notes: \_\_\_\_\_  
 \_\_\_\_\_  
 \_\_\_\_\_

**VALIDATION FINDINGS WORKSHEET**  
**Sample Specific Analysis Reference**

All circled methods are applicable to each sample.

Sample ID	Matrix	Parameter
1 → 29	W	pH (TDS) CI F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> (ClO <sub>4</sub> )
QC 30,31	↓	pH (TDS) CI F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
↓ 32,33	↓	pH TDS CI F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> (ClO <sub>4</sub> )
		pH TDS CI F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
		pH TDS CI F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
		pH TDS CI F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
		pH TDS CI F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
		pH TDS CI F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
		pH TDS CI F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
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		pH TDS CI F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
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		pH TDS CI F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
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Comments:

VALIDATION FINDINGS WORKSHEET  
Technical Holding Times

All circled dates have exceeded the technical holding time.  
Y N N/A Were all samples preserved as applicable to each method?  
Y N N/A Were all cooler temperatures within validation criteria?

Method:	314.0					
Parameters:	C104					
Technical holding time:	28 days					

Sample ID	Sampling date	Analysis date	Analysis date	Analysis date	Analysis date	Analysis date	Qualifier
26	8-11-10	9-9-10	(29 days)				J-UJA h
		(reanalysis)					

LDC #: 74203 J6  
SDG #:           

### VALIDATION FINDINGS WORKSHEET Laboratory Control Samples (LCS)

Page: 1 of 1  
Reviewer: MCF  
2nd Reviewer:    

METHOD: Inorganics, Method See cover

Please see qualifications below for all questions answered "N". Not applicable questions are identified as "N/A".

Y N N/A Was a laboratory control sample (LCS) analyzed for each matrix in this SDG?

Y N/A Were the LCS percent recoveries (%R) and relative percent difference (RPD) within the QC limits?

LEVEL IV ONLY:

Y N N/A Were recalculated results acceptable? See Level IV Recalculation Worksheet for recalculations.

#	LCS/LCSD ID	Matrix	Analyte	LCS %R (limits)	LCSD %R (limits)	RPD (limits)	Associated Samples	Qualifications
1	LCS/LCSD	Water	ClO4	79 (85-115)	26 (≤ 20)	4, 11 → 24, 28, 29		* No Qual F/MS/P (e, ed)

Comments: ~~\* < LCS~~ 93 see in limit

**Laboratory Data Consultants, Inc.  
Data Validation Report**

**Project/Site Name:** 2010 Annual Remedial Performance Sampling  
**Collection Date:** August 23, 2010  
**LDC Report Date:** December 27, 2010  
**Matrix:** Water  
**Parameters:** Wet Chemistry  
**Validation Level:** Stage 2A  
**Laboratory:** MWH Laboratories, Inc.  
**Sample Delivery Group (SDG):** 341684

**Sample Identification**

PC-144  
PC-145

## Introduction

This data review covers 2 water samples listed on the cover sheet including dilutions and reanalysis as applicable. The analyses were per EPA Method 314.0 for Perchlorate, and EPA Method 160.1 and Standard Method 2540C for Total Dissolved Solids.

This review follows a modified outline of the USEPA Contract Laboratory Program National Functional Guidelines for Inorganic Data Review (October 2004) and the EPA Region 9 Superfund Data Evaluation/Validation Guidance, NDEP guidance (May 2006).

A qualification summary table is provided at the end of this report if data has been qualified. Flags are classified as P (protocol) or A (advisory) to indicate whether the flag is due to a laboratory deviation from a specified protocol or is of technical advisory nature.

Raw data were not reviewed for this SDG. The review was based on QC data.

The following are definitions of the data qualifiers:

- J+ Data are qualified as estimated, with a high bias likely to occur. False positives or false negatives are unlikely to have been reported.
- J- Data are qualified as estimated, with a low bias likely to occur. False positives or false negatives are unlikely to have been reported.
- J Data are qualified as estimated; it is not possible to assess the direction of the potential bias. False positives or false negatives are unlikely to have been reported.
- U Indicates the compound or analyte was analyzed for but not detected at or above the stated limit.
- R Data are qualified as rejected. There is a significant potential for the reporting of false negatives or false positives.
- UJ Indicates the compound or analyte was analyzed for but not detected. The sample detection limit is an estimated value.
- A Indicates the finding is based upon technical validation criteria.
- P Indicates the finding is related to a protocol/contractual deviation.
- None Indicates the data was not significantly impacted by the finding, therefore qualification was not required.

## **I. Technical Holding Times**

All technical holding time requirements were met.

The chain-of-custodies were reviewed for documentation of cooler temperatures. All cooler temperatures met validation criteria.

## **II. Calibration**

### **a. Initial Calibration**

Initial calibration data were not reviewed for Stage 2A.

### **b. Calibration Verification**

Calibration verification data were not reviewed for Stage 2A.

## **III. Blanks**

Method blanks were reviewed for each matrix as applicable. No contaminant concentrations were found in the preparation blanks.

No field blanks were identified in this SDG.

## **IV. Matrix Spike/Matrix Spike Duplicates**

The laboratory has indicated that there were no matrix spike (MS) analyses specified for the samples in this SDG, and therefore matrix spike analyses were not performed for this SDG.

## **V. Duplicates**

The laboratory has indicated that there were no duplicate (DUP) analyses specified for the samples in this SDG, and therefore duplicate analyses were not performed for this SDG.

## **VI. Laboratory Control Samples**

Laboratory control samples were reviewed for each matrix as applicable. Percent recoveries (%R) and relative percent differences (RPD) were within QC limits.

## **VII. Sample Result Verification**

Raw data were not reviewed for this SDG.

## **VIII. Overall Assessment of Data**

Data flags are summarized at the end of this report if data has been qualified.

## **IX. Field Duplicates**

No field duplicates were identified in this SDG.



**2010 Annual Remedial Performance Sampling  
Wet Chemistry - Data Qualification Summary - SDG 341684**

No Sample Data Qualified in this SDG

**2010 Annual Remedial Performance Sampling  
Wet Chemistry - Laboratory Blank Data Qualification Summary - SDG 341684**

No Sample Data Qualified in this SDG

**2010 Annual Remedial Performance Sampling  
Wet Chemistry - Field Blank Data Qualification Summary - SDG 341684**

No Sample Data Qualified in this SDG

LDC #: 24203K6  
 SDG #: 341684  
 Laboratory: MWH Laboratories

**Tronox Northgate Henderson**  
**VALIDATION COMPLETENESS WORKSHEET**

Stage 2 ~~B~~ A  
 qm

Date: 12-2-10  
 Page: 1 of 1  
 Reviewer: MG  
 2nd Reviewer: [Signature]

**METHOD: (Analyte)** Perchlorate (EPA Method 314.0), TDS (EPA Method 160.1/SM2540C)

The samples listed below were reviewed for each of the following validation areas. Validation findings are noted in attached validation findings worksheets.

	Validation Area		Comments
I.	Technical holding times	A	Sampling dates: <u>8-23-10</u>
IIa.	Initial calibration	N	
IIb.	Calibration verification	N	
III.	Blanks	A	
IV	Matrix Spike/Matrix Spike Duplicates	N	<u>client specified</u>
V	Duplicates	N	" "
VI.	Laboratory control samples	A	<u>LCS/LCSD</u>
VII.	Sample result verification	N	
VIII.	Overall assessment of data	A	
IX.	Field duplicates	N	
X	Field blanks	N	

Note: A = Acceptable                      ND = No compounds detected                      D = Duplicate  
 N = Not provided/applicable                      R = Rinsate                      TB = Trip blank  
 SW = See worksheet                      FB = Field blank                      EB = Equipment blank

Validated Samples:  
all water

1	PC-144	11		21		31	
2	PC-145	12		22		32	
3	<u>PBW</u>	13		23		33	
4		14		24		34	
5		15		25		35	
6		16		26		36	
7		17		27		37	
8		18		28		38	
9		19		29		39	
10		20		30		40	

Notes: \_\_\_\_\_  
 \_\_\_\_\_  
 \_\_\_\_\_

VALIDATION FINDINGS WORKSHEET  
Sample Specific Analysis Reference

All circled methods are applicable to each sample.

Sample ID	Matrix	Parameter
1, 2	W	pH TDS CI F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> CIO <sub>4</sub>
		pH TDS CI F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> CIO <sub>4</sub>
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Comments: \_\_\_\_\_

**Laboratory Data Consultants, Inc.  
Data Validation Report**

**Project/Site Name:** 2010 Annual Remedial Performance Sampling  
**Collection Date:** September 8, 2010  
**LDC Report Date:** December 19, 2010  
**Matrix:** Water  
**Parameters:** Wet Chemistry  
**Validation Level:** Stage 2A  
**Laboratory:** MWH Laboratories, Inc.  
**Sample Delivery Group (SDG):** 343130

**Sample Identification**

ART-1  
ART-2  
ART-3  
ART-4  
ART-6  
ART-7  
ART-8  
PC-99R2/R3  
PC-115R  
PC-116R  
SF-1  
PC-117  
PC-118  
PC-119  
PC-120  
PC-121  
PC-133  
ART-9

## Introduction

This data review covers 18 water samples listed on the cover sheet including dilutions and reanalysis as applicable. The analyses were per EPA Method 314.0 for Perchlorate, and EPA Method 160.1 and Standard Method 2540C for Total Dissolved Solids.

This review follows a modified outline of the USEPA Contract Laboratory Program National Functional Guidelines for Inorganic Data Review (October 2004) and the EPA Region 9 Superfund Data Evaluation/Validation Guidance, NDEP guidance (May 2006).

A qualification summary table is provided at the end of this report if data has been qualified. Flags are classified as P (protocol) or A (advisory) to indicate whether the flag is due to a laboratory deviation from a specified protocol or is of technical advisory nature.

Raw data were not reviewed for this SDG. The review was based on QC data.

The following are definitions of the data qualifiers:

- J+ Data are qualified as estimated, with a high bias likely to occur. False positives or false negatives are unlikely to have been reported.
- J- Data are qualified as estimated, with a low bias likely to occur. False positives or false negatives are unlikely to have been reported.
- J Data are qualified as estimated; it is not possible to assess the direction of the potential bias. False positives or false negatives are unlikely to have been reported.
- U Indicates the compound or analyte was analyzed for but not detected at or above the stated limit.
- R Data are qualified as rejected. There is a significant potential for the reporting of false negatives or false positives.
- UJ Indicates the compound or analyte was analyzed for but not detected. The sample detection limit is an estimated value.
- A Indicates the finding is based upon technical validation criteria.
- P Indicates the finding is related to a protocol/contractual deviation.
- None Indicates the data was not significantly impacted by the finding, therefore qualification was not required.

## **I. Technical Holding Times**

All technical holding time requirements were met.

The chain-of-custodies were reviewed for documentation of cooler temperatures. All cooler temperatures met validation criteria.

## **II. Calibration**

### **a. Initial Calibration**

Initial calibration data were not reviewed for Stage 2A.

### **b. Calibration Verification**

Calibration verification data were not reviewed for Stage 2A.

## **III. Blanks**

Method blanks were reviewed for each matrix as applicable. No contaminant concentrations were found in the preparation blanks.

No field blanks were identified in this SDG.

## **IV. Matrix Spike/Matrix Spike Duplicates**

The laboratory has indicated that there were no matrix spike (MS) analyses specified for the samples in this SDG, and therefore matrix spike analyses were not performed for this SDG.

## **V. Duplicates**

The laboratory has indicated that there were no duplicate (DUP) analyses specified for the samples in this SDG, and therefore duplicate analyses were not performed for this SDG.

## **VI. Laboratory Control Samples**

Laboratory control samples were reviewed for each matrix as applicable. Percent recoveries (%R) and relative percent differences (RPD) were within QC limits.

## **VII. Sample Result Verification**

Raw data were not reviewed for this SDG.

## **VIII. Overall Assessment of Data**

Data flags are summarized at the end of this report if data has been qualified.

## **IX. Field Duplicates**

No field duplicates were identified in this SDG.

**2010 Annual Remedial Performance Sampling  
Wet Chemistry - Data Qualification Summary - SDG 343130**

No Sample Data Qualified in this SDG

**2010 Annual Remedial Performance Sampling  
Wet Chemistry - Laboratory Blank Data Qualification Summary - SDG 343130**

No Sample Data Qualified in this SDG

**2010 Annual Remedial Performance Sampling  
Wet Chemistry - Field Blank Data Qualification Summary - SDG 343130**

No Sample Data Qualified in this SDG



LDC #: 24203L6

**Tronox Northgate Henderson  
VALIDATION COMPLETENESS WORKSHEET**

Date: 12-2-10

SDG #: 343130

Stage ~~2~~ **A**

Page: 1 of 1

Laboratory: MWH Laboratories

*AM*

Reviewer: MG

2nd Reviewer: ✓

**METHOD: (Analyte)** Perchlorate (EPA Method 314.0), TDS (EPA Method 160.1/SM2540C)

The samples listed below were reviewed for each of the following validation areas. Validation findings are noted in attached validation findings worksheets.

	Validation Area		Comments
I.	Technical holding times	A	Sampling dates: <u>9-8-10</u>
IIa.	Initial calibration	N	
IIb.	Calibration verification	N	
III.	Blanks	A	
IV	Matrix Spike/Matrix Spike Duplicates	N	<u>client specified</u>
V	Duplicates	N	<u>" "</u>
VI.	Laboratory control samples	A	<u>LCS/LCSD</u>
VII.	Sample result verification	N	
VIII.	Overall assessment of data	A	
IX.	Field duplicates	N	
X	Field blanks	N	

Note: A = Acceptable  
N = Not provided/applicable  
SW = See worksheet

ND = No compounds detected  
R = Rinsate  
FB = Field blank

D = Duplicate  
TB = Trip blank  
EB = Equipment blank

Validated Samples:  
all water

1	ART-1	11	SF-1	21		31	
2	ART-2	12	PC-117	22		32	
3	ART-3	13	PC-118	23		33	
4	ART-4	14	PC-119	24		34	
5	ART-6	15	PC-120	25		35	
6	ART-7	16	PC-121	26		36	
7	ART-8	17	PC-133	27		37	
8	PC-99R2/R3	18	ART-9	28		38	
9	PC-115R	19	<u>PBW</u>	29		39	
10	PC-116R	20		30		40	

Notes: \_\_\_\_\_  
\_\_\_\_\_  
\_\_\_\_\_

LDC #: 24203L6

**VALIDATION FINDINGS WORKSHEET**  
**Sample Specific Analysis Reference**

Page: 1 of 1  
Reviewer: MG  
2nd reviewer: W

All circled methods are applicable to each sample.

Sample ID	Matrix	Parameter
1 → 18	W	pH TDS <u>Cl</u> F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> <u>ClO<sub>4</sub></u>
		pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
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		pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
		pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
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Comments: \_\_\_\_\_

## Laboratory Data Consultants, Inc. Data Validation Report

**Project/Site Name:** 2010 Annual Remedial Performance Sampling  
**Collection Date:** September 13 through September 16, 2010  
**LDC Report Date:** December 27, 2010  
**Matrix:** Water  
**Parameters:** Wet Chemistry  
**Validation Level:** Stage 2A  
**Laboratory:** MWH Laboratories, Inc.  
**Sample Delivery Group (SDG):** 343913

### Sample Identification

M-83	PC-53
M-87	PC-103
PC-98R	MW-K5
PC-86	PC-91
PC-90	PC-97
PC-56	PC-18
PC-58	PC-55
PC-59	PC-101R
PC-60	M-83DUP
PC-62	PC-68DUP
PC-68	PC-53DUP
PC-122	
MW-K4	
ARP-1	
ARP-2A	
ARP-3A	
ARP-4A	
ARP-5A	
ARP-6B	
ARP-7	

## Introduction

This data review covers 31 water samples listed on the cover sheet including dilutions and reanalysis as applicable. The analyses were per EPA Method 314.0 for Perchlorate, and EPA Method 160.1 and Standard Method 2540C for Total Dissolved Solids.

This review follows a modified outline of the USEPA Contract Laboratory Program National Functional Guidelines for Inorganic Data Review (October 2004) and the EPA Region 9 Superfund Data Evaluation/Validation Guidance, NDEP guidance (May 2006).

A qualification summary table is provided at the end of this report if data has been qualified. Flags are classified as P (protocol) or A (advisory) to indicate whether the flag is due to a laboratory deviation from a specified protocol or is of technical advisory nature.

Raw data were not reviewed for this SDG. The review was based on QC data.

The following are definitions of the data qualifiers:

- J+ Data are qualified as estimated, with a high bias likely to occur. False positives or false negatives are unlikely to have been reported.
- J- Data are qualified as estimated, with a low bias likely to occur. False positives or false negatives are unlikely to have been reported.
- J Data are qualified as estimated; it is not possible to assess the direction of the potential bias. False positives or false negatives are unlikely to have been reported.
- U Indicates the compound or analyte was analyzed for but not detected at or above the stated limit.
- R Data are qualified as rejected. There is a significant potential for the reporting of false negatives or false positives.
- UJ Indicates the compound or analyte was analyzed for but not detected. The sample detection limit is an estimated value.
- A Indicates the finding is based upon technical validation criteria.
- P Indicates the finding is related to a protocol/contractual deviation.
- None Indicates the data was not significantly impacted by the finding, therefore qualification was not required.

## **I. Technical Holding Times**

All technical holding time requirements were met.

The chain-of-custodies were reviewed for documentation of cooler temperatures. All cooler temperatures met validation criteria.

## **II. Calibration**

### **a. Initial Calibration**

Initial calibration data were not reviewed for Stage 2A.

### **b. Calibration Verification**

Calibration verification data were not reviewed for Stage 2A.

## **III. Blanks**

Method blanks were reviewed for each matrix as applicable. No contaminant concentrations were found in the preparation blanks.

No field blanks were identified in this SDG.

## **IV. Matrix Spike/Matrix Spike Duplicates**

The laboratory has indicated that there were no matrix spike (MS) analyses specified for the samples in this SDG, and therefore matrix spike analyses were not performed for this SDG.

## **V. Duplicates**

Duplicate (DUP) sample analyses were reviewed for each matrix as applicable. Results were within QC limits.

## **VI. Laboratory Control Samples**

Laboratory control samples were reviewed for each matrix as applicable. Percent recoveries (%R) and relative percent differences (RPD) were within QC limits with the following exceptions:

LCS ID (Associated Samples)	Analyte	LCS %R (Limits)	LCSD %R (Limits)	RPD (Limits)	Flag	A or P
LCS/LCSD (M-87 PC-98R PC-59 PC-60 PC-62)	Perchlorate	-	83 (84-115)	23 ( $\leq 20$ )	J (all detects) UJ (all non-detects)	P

## VII. Sample Result Verification

Raw data were not reviewed for this SDG.

## VIII. Overall Assessment of Data

Data flags are summarized at the end of this report if data has been qualified.

## IX. Field Duplicates

No field duplicates were identified in this SDG.

**2010 Annual Remedial Performance Sampling  
Wet Chemistry - Data Qualification Summary - SDG 343913**

SDG	Sample	Analyte	Flag	A or P	Reason
343913	M-87 PC-98R PC-59 PC-60 PC-62	Perchlorate	J (all detects) UJ (all non-detects)	P	Laboratory control samples (%R) (RPD)

**2010 Annual Remedial Performance Sampling  
Wet Chemistry - Laboratory Blank Data Qualification Summary - SDG 343913**

No Sample Data Qualified in this SDG

**2010 Annual Remedial Performance Sampling  
Wet Chemistry - Field Blank Data Qualification Summary - SDG 343913**

No Sample Data Qualified in this SDG

**Tronox Northgate Henderson**

**VALIDATION COMPLETENESS WORKSHEET**

LDC #: 24203M6  
 SDG #: 343913  
 Laboratory: MWH Laboratories

Stage ~~2B~~ **A**  
*9/14*

Date: 12-2-10  
 Page: 1 of 1  
 Reviewer: MG  
 2nd Reviewer: [Signature]

**METHOD: (Analyte)** Perchlorate (EPA Method 314.0), TDS (EPA Method 160.1/SM2540C)

The samples listed below were reviewed for each of the following validation areas. Validation findings are noted in attached validation findings worksheets.

	Validation Area		Comments
I.	Technical holding times	A	Sampling dates: <u>9-13-10 through 9-16-10</u>
IIa.	Initial calibration	N	
IIb.	Calibration verification	N	
III.	Blanks	A	
IV	Matrix Spike/Matrix Spike Duplicates	N	<u>client specified</u>
V	Duplicates	A	<u>DUP</u>
VI.	Laboratory control samples	SW	<u>LCS/LCSD</u>
VII.	Sample result verification	N	
VIII.	Overall assessment of data	A	
IX.	Field duplicates	N	
X	Field blanks	N	

Note: A = Acceptable                      ND = No compounds detected                      D = Duplicate  
 N = Not provided/applicable                      R = Rinsate                      TB = Trip blank  
 SW = See worksheet                      FB = Field blank                      EB = Equipment blank

Validated Samples: all water

1	M-83	11	PC-68	21	PC-53	31	PC-53DUP
2	M-87	12	PC-122	22	PC-103	32	
3	PC-98R	13	MW-K4	23	MW-K5	33	
4	PC-86	14	ARP-1	24	PC-91	34	
5	PC-90	15	ARP-2A	25	PC-97	35	<u>PBW1</u>
6	PC-56	16	ARP-3A	26	PC-18	36	<u>PBW2</u>
7	PC-58	17	ARP-4A	27	PC-55	37	<u>PBW3</u>
8	PC-59	18	ARP-5A	28	PC-101R	38	<u>PBW4</u>
9	PC-60	19	<sup>6B</sup> <del>ARP-86</del> <i>9/14</i>	29	M-83DUP	39	<u>PBW5</u>
10	PC-62	20	ARP-7	30	PC-68DUP	40	<u>PBW6</u>

Notes: \_\_\_\_\_  
 \_\_\_\_\_  
 \_\_\_\_\_





LDC #: 24203M6

SDG #:         

**VALIDATION FINDINGS WORKSHEET**  
**Laboratory Control Samples (LCS)**

Page: 1 of 1  
Reviewer: MG  
2nd Reviewer:         

METHOD: Inorganics, Method          see cover

Please see qualifications below for all questions answered "N". Not applicable questions are identified as "N/A".

(Y) N N/A Was a laboratory control sample (LCS) analyzed for each matrix in this SDG?

Y (N) N/A Were the LCS percent recoveries (%R) and relative percent difference (RPD) within the QC limits?

**LEVEL IV ONLY:**

Y N N/A Were recalculated results acceptable? See Level IV Recalculation Worksheet for recalculations.

#	LCS/LCSD ID	Matrix	Analyte	LCS %R (limits)	LCSD %R (limits)	RPD (limits)	Associated Samples	Qualifications
1	LCS/LCSD	Water	ClO4		83 (85-115)	23 (≤ 20)	2, 3, 8 → 10	0* No Quant-1/10/14 (S, LD)

Comments: \* LCS %a rec in limit e

**Laboratory Data Consultants, Inc.  
Data Validation Report**

**Project/Site Name:** 2010 Annual Remedial Performance Sampling  
**Collection Date:** August 1 through August 2, 2010  
**LDC Report Date:** December 23, 2010  
**Matrix:** Water  
**Parameters:** Wet Chemistry  
**Validation Level:** Stage 2A  
**Laboratory:** MWH Laboratories, Inc.  
**Sample Delivery Group (SDG):** 339791

**Sample Identification**

I-O	PC-132
I-P	M-96
I-H	PC-54
I-U	PC-37
I-T	PC-71
I-G	PC-72
I-F	PC-73
I-N	M-23
I-E	VD080210
I-M	FB080210V
I-D	I-OMS
PC-123	I-OMSD
PC-124	I-ODUP
PC-125	I-DMS
PC-126	I-DMSD
PC-127	PC-125DUP
PC-128	PC-54DUP
PC-129	FB080210VMS
PC-130	FB080210VMDS
PC-131	

## Introduction

This data review covers 39 water samples listed on the cover sheet including dilutions and reanalysis as applicable. The analyses were per EPA SW 846 Method 7196A for Hexavalent Chromium, EPA Method 314.0 for Perchlorate, and EPA Method 160.1 and Standard Method 2540C for Total Dissolved Solids.

This review follows a modified outline of the USEPA Contract Laboratory Program National Functional Guidelines for Inorganic Data Review (October 2004) and the EPA Region 9 Superfund Data Evaluation/Validation Guidance, NDEP guidance (May 2006).

A qualification summary table is provided at the end of this report if data has been qualified. Flags are classified as P (protocol) or A (advisory) to indicate whether the flag is due to a laboratory deviation from a specified protocol or is of technical advisory nature.

Raw data were not reviewed for this SDG. The review was based on QC data.

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- A Indicates the finding is based upon technical validation criteria.
- P Indicates the finding is related to a protocol/contractual deviation.
- None Indicates the data was not significantly impacted by the finding, therefore qualification was not required.

## I. Technical Holding Times

All technical holding time requirements were met with the following exceptions:

Sample	Analyte	Total Time From Sample Collection Until Analysis	Required Holding Time From Sample Collection Until Analysis	Flag	A or P
I-O I-P I-H I-U I-T I-G I-F I-N I-E I-M I-OMS I-OMSD	Perchlorate	70 days	28 days	J- (all detects) R (all non-detects)	P
I-D PC-54 PC-37 PC-71 PC-72 PC-73 I-DMS I-DMSD	Perchlorate	71 days	28 days	J- (all detects) R (all non-detects)	P
PC-123 PC-124 PC-125 PC-126 PC-127 PC-128 PC-129 PC-130 PC-132 M-96 M-23 VD080210 FB080210V	Perchlorate	72 days	28 days	J- (all detects) R (all non-detects)	P
FB080210V FB080210VMS FB080210VMSD	Hexavalent chromium	34.75 hours	24 hours	J- (all detects) UJ (all nondetects)	P

The chain-of-custodies were reviewed for documentation of cooler temperatures. All cooler temperatures met validation criteria.

## II. Calibration

### a. Initial Calibration

Initial calibration data were not reviewed for Stage 2A.

## b. Calibration Verification

Calibration verification data were not reviewed for Stage 2A.

## III. Blanks

Method blanks were reviewed for each matrix as applicable. No contaminant concentrations were found in the preparation blanks.

Sample FB080210V was identified as a field blank. No contaminant concentrations were found in this blank with the following exceptions:

Field Blank ID	Sampling Date	Analyte	Concentration	Associated Samples
FB080210V	8/2/10	Hexavalent chromium	0.0060 mg/L	No associated sample in this SDG
FB080210V	8/2/10	Perchlorate	1.63 ug/L	I-O I-P I-H I-U I-T I-G I-F I-N I-E I-M I-D PC-54 PC-37 PC-71 PC-72 PC-73 M-23 VD080210

Sample concentrations were compared to concentrations detected in the field blanks as required by the QAPP. No sample data was qualified.

## IV. Matrix Spike/Matrix Spike Duplicates

Matrix spike (MS) and matrix spike duplicate (MSD) samples were reviewed for each matrix as applicable. Percent recoveries (%R) and relative percent differences (RPD) were within QC limits.

## V. Duplicates

Duplicate (DUP) sample analyses were reviewed for each matrix as applicable. Results were within QC limits.

## VI. Laboratory Control Samples

Laboratory control samples were reviewed for each matrix as applicable. Percent recoveries (%R) and relative percent differences (RPD) were within QC limits.

## VII. Sample Result Verification

Raw data were not reviewed for this SDG.

## VIII. Overall Assessment of Data

Data flags are summarized at the end of this report if data has been qualified.

## IX. Field Duplicates

Samples M-23 and VD080210 were identified as field duplicates. No contaminant concentrations were detected in any of the samples with the following exceptions:

Analyte	Concentration		RPD (Limits)	Difference (Limits)	Flags	A or P
	M-23	VD080210				
Perchlorate	310000 ug/L	310000 ug/L	0 (≤30)	-	-	-
Total dissolved solids	4600 mg/L	4600 mg/L	0 (≤30)	-	-	-

**2010 Annual Remedial Performance Sampling  
Wet Chemistry - Data Qualification Summary - SDG 339791**

SDG	Sample	Analyte	Flag	A or P	Reason
339791	I-O I-P I-H I-U I-T I-G I-F I-N I-E I-M I-D PC-54 PC-37 PC-71 PC-72 PC-73 PC-123 PC-124 PC-125 PC-126 PC-127 PC-128 PC-129 PC-130 PC-132 M-96 M-23 VD080210 FB080210V	Perchlorate	J- (all detects) R (all non-detects)	P	Technical holding times
339791	FB080210V	Hexavalent chromium	J- (all detects) UJ (all nondetects)	P	Technical holding times

**2010 Annual Remedial Performance Sampling  
Wet Chemistry - Laboratory Blank Data Qualification Summary - SDG 339791**

No Sample Data Qualified in this SDG

**2010 Annual Remedial Performance Sampling  
Wet Chemistry - Field Blank Data Qualification Summary - SDG 339791**

No Sample Data Qualified in this SDG



Tronox Northgate Henderson

VALIDATION COMPLETENESS WORKSHEET

LDC #: 24203N6  
 SDG #: 339791  
 Laboratory: MWH Laboratories

Stage 2B A

SMY

Date: 12-15-10  
 Page: 1 of 1  
 Reviewer: MG  
 2nd Reviewer: W

METHOD: (Analyte) Hexavalent Chromium (EPA SW846 Method 7196A), Perchlorate (EPA Method 314.0), TDS (EPA Method 160.1/SM2540C)

The samples listed below were reviewed for each of the following validation areas. Validation findings are noted in attached validation findings worksheets.

	Validation Area		Comments
I.	Technical holding times	SW	Sampling dates: 8-1-10 through 8-2-10
IIa.	Initial calibration	N	
IIb.	Calibration verification	N	
III.	Blanks	A	
IV	Matrix Spike/Matrix Spike Duplicates	A	MS/MSD
V	Duplicates	A	DUP. (SDG: 339977)
VI.	Laboratory control samples	A	LCS/LCSD
VII.	Sample result verification	N	
VIII.	Overall assessment of data	A	
IX.	Field duplicates	SW	D = 28 + 29
X	Field blanks	SW	FB = 30

Note: A = Acceptable  
 N = Not provided/applicable  
 SW = See worksheet  
 ND = No compounds detected  
 R = Rinsate  
 FB = Field blank  
 D = Duplicate  
 TB = Trip blank  
 EB = Equipment blank

Validated Samples:  
 all water

1	I-O	11	I-D	21	PC-132	31	I-OMS
2	I-P	12	PC-123	22	M-96	32	I-OMSD
3	I-H	13	PC-124	23	PC-54	33	I-ODUP
4	I-U	14	PC-125	24	PC-37	34	I-DMS
5	I-T	15	PC-126	25	PC-71	35	I-DMSD
6	I-G	16	PC-127	26	PC-72	36	PC-125DUP
7	I-F	17	PC-128	27	PC-73	37	PC-54DUP
8	I-N	18	PC-129	28	M-23	38	FB080210VMS
9	I-E	19	PC-130	29	VD080210	39	FB080210VMSD
10	I-M	20	PC-131	30	FB080210V	40	

Notes: PBW1  
 PBW2  
 PBW3

**VALIDATION FINDINGS WORKSHEET**  
**Sample Specific Analysis Reference**

All circled methods are applicable to each sample.

Sample ID	Matrix	Parameter
1 → 29	W	pH (TDS) Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
30	↓	pH (TDS) Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
OC 31, 32, 34, 35		pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
33, 36, 37	↓	pH (TDS) Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
38, 39		pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
		pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
		pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
		pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
		pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
		pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
		pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
		pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
		pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
		pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
		pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
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		pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
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		pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
		pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
		pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
		pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
		pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
		pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
		pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
		pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
		pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
		pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>

Comments: \_\_\_\_\_

### VALIDATION FINDINGS WORKSHEET Technical Holding Times

All circled dates have exceeded the technical holding time.

N N/A Were all samples preserved as applicable to each method?

N N/A Were all cooler temperatures within validation criteria?

Method:	314.0		7196A				
Parameters:	C104		Cr VI				
Technical holding time:	28 days		24 hr				
Sample ID	Sampling date	Analysis date	Analysis date	Analysis date	Analysis date	Analysis date	Qualifier
1 → 10, 31, 32	8-2-10	10-11-10	(70 days)				J-R/P
11, 23 → 27, 34, 35	↓	10-12-10	(71 days)				↓
12 → 19, 21, 22	8-1-10	↓	(72 days)				↓
28 → 30	8-2-10	10-13-10	( ↓ )				↓
30	<sup>0241</sup> 8-2-10			<sup>13.33</sup> 8-3-10	(34.75 hr)		J-UJ/P
38	↓			↓	↓		↓
39	↓			↓	↓		↓

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**VALIDATION FINDINGS WORKSHEET**  
Field Blanks

**METHOD:** Inorganics, EPA Method See Cover  
 N N/A Were field blanks identified in this SDG?  
 N N/A Were target analytes detected in the field blanks?  
**Blank units:** ug/L Associated sample units: ug/L  
**Sampling date:** 8/2/10 Soil factor applied NA  
**Field blank type:** (circle one) Field Blank / Rinsate / Other:

Associated Samples: CIO4: 1-11, 23-29 (>RL), Cr VI: none

Analyte	Blank ID	Action Limit	Sample Identification							
	30	No Qual's.								
Cr VI (mg/L)	0.0060	0.06								
CIO4	1.63									

CIRCLED RESULTS WERE NOT QUALIFIED. ALL RESULTS NOT CIRCLED WERE QUALIFIED BY THE FOLLOWING STATEMENT:  
 Samples with analyte concentrations within five times the associated field blank concentration are listed above, these sample results were qualified as not detected, "U".

LDC#: 24203N6  
SDG#: See Cover

**VALIDATION FINDINGS WORKSHEET**  
**Field Duplicates**

Page: 1 of 1  
Reviewer: MC  
2nd Reviewer: [Signature]

Inorganics, Method See Cover

- Y N NA Were field duplicate pairs identified in this SDG?
- Y N NA Were target analytes detected in the field duplicate pairs?

Analyte	Concentration (mg/L)		RPD ( $\leq 30$ )	Difference	Limits	Qualification (Parent only)
	28	29				
Perchlorate (ug/L)	310000	310000	0			
TDS	4600	4600	0			

V:\FIELD DUPLICATES\FD\_inorganic\24203N6.wpd

**Laboratory Data Consultants, Inc.  
Data Validation Report**

**Project/Site Name:** 2010 Annual Remedial Performance Sampling  
**Collection Date:** August 2 through August 3, 2010  
**LDC Report Date:** December 27, 2010  
**Matrix:** Water  
**Parameters:** Wet Chemistry  
**Validation Level:** Stage 2A  
**Laboratory:** MWH Laboratories, Inc.  
**Sample Delivery Group (SDG):** 339977

**Sample Identification**

I-Q	EB080310VMS
I-C	EB080310VMDS
I-S	
I-L	
I-R	
I-B	
I-AR	
I-AB	
I-AA	
M-131	
M-57A	
M-79	
M-69	
M-135	
VD-080310	
EB080310V	
I-CMS	
I-CMSD	
I-CDUP	
M-79DUP	

## Introduction

This data review covers 22 water samples listed on the cover sheet including dilutions and reanalysis as applicable. The analyses were per EPA SW 846 Method 7196A for Hexavalent Chromium, EPA Method 314.0 for Perchlorate, and EPA Method 160.1 and Standard Method 2540C for Total Dissolved Solids.

This review follows a modified outline of the USEPA Contract Laboratory Program National Functional Guidelines for Inorganic Data Review (October 2004) and the EPA Region 9 Superfund Data Evaluation/Validation Guidance, NDEP guidance (May 2006).

A qualification summary table is provided at the end of this report if data has been qualified. Flags are classified as P (protocol) or A (advisory) to indicate whether the flag is due to a laboratory deviation from a specified protocol or is of technical advisory nature.

Raw data were not reviewed for this SDG. The review was based on QC data.

The following are definitions of the data qualifiers:

- J+ Data are qualified as estimated, with a high bias likely to occur. False positives or false negatives are unlikely to have been reported.
- J- Data are qualified as estimated, with a low bias likely to occur. False positives or false negatives are unlikely to have been reported.
- J Data are qualified as estimated; it is not possible to assess the direction of the potential bias. False positives or false negatives are unlikely to have been reported.
- U Indicates the compound or analyte was analyzed for but not detected at or above the stated limit.
- R Data are qualified as rejected. There is a significant potential for the reporting of false negatives or false positives.
- UJ Indicates the compound or analyte was analyzed for but not detected. The sample detection limit is an estimated value.
- A Indicates the finding is based upon technical validation criteria.
- P Indicates the finding is related to a protocol/contractual deviation.
- None Indicates the data was not significantly impacted by the finding, therefore qualification was not required.

## I. Technical Holding Times

All technical holding time requirements were met with the following exceptions:

Sample	Analyte	Total Time From Sample Collection Until Analysis	Required Holding Time From Sample Collection Until Analysis	Flag	A or P
I-Q I-C I-S I-L I-R I-B I-AR I-AB I-CMS I-CMSD M-131	Perchlorate	72 days	28 days	J- (all detects) R (all non-detects)	P
I-AA	Perchlorate	73 days	28 days	J- (all detects) R (all non-detects)	P
M-57A M-79 M-69 M-135 VD-080310	Perchlorate	71 days	28 days	J- (all detects) R (all non-detects)	P
EB080310V EB080310VMS EB080310VMSD	Hexavalent chromium	61 hours	24 hours	J- (all detects) R (all non-detects)	P

The chain-of-custodies were reviewed for documentation of cooler temperatures. All cooler temperatures met validation criteria.

## II. Calibration

### a. Initial Calibration

Initial calibration data were not reviewed for Stage 2A.

### b. Calibration Verification

Calibration verification data were not reviewed for Stage 2A.

## III. Blanks

Method blanks were reviewed for each matrix as applicable. No contaminant concentrations were found in the preparation blanks.

Sample EB080310V was identified as an equipment blank. No contaminant concentrations were found in this blank with the following exceptions:



Equipment Blank ID	Sampling Date	Analyte	Concentration	Associated Samples
EB080310V	8/3/10	Perchlorate	8.5 ug/L	M-131 M-57A M-79 M-69 M-135 VD-080310

Sample FB080210V (from SDG 339791) was identified as a field blank. No contaminant concentrations were found in this blank with the following exceptions:

Field Blank ID	Sampling Date	Analyte	Concentration	Associated Samples
FB080210V	8/2/10	Hexavalent chromium	0.0060 mg/L	No associated sample in this SDG
FB080210V	8/2/10	Perchlorate	1.63 ug/L	I-Q I-C I-S I-L I-R I-B I-AR I-AB I-AA

Sample concentrations were compared to concentrations detected in the field blanks as required by the QAPP. No sample data was qualified.

#### IV. Matrix Spike/Matrix Spike Duplicates

Matrix spike (MS) and matrix spike duplicate (MSD) samples were reviewed for each matrix as applicable. Percent recoveries (%R) and relative percent differences (RPD) were within QC limits.

#### V. Duplicates

Duplicate (DUP) sample analyses were reviewed for each matrix as applicable. Results were within QC limits.

#### VI. Laboratory Control Samples

Laboratory control samples were reviewed for each matrix as applicable. Percent recoveries (%R) and relative percent differences (RPD) were within QC limits.

#### VII. Sample Result Verification

Raw data were not reviewed for this SDG.

### VIII. Overall Assessment of Data

Data flags are summarized at the end of this report if data has been qualified.

### IX. Field Duplicates

Samples M-131 and VD-080310 were identified as field duplicates. No contaminant concentrations were detected in any of the samples with the following exceptions:

Analyte	Concentration		RPD (Limits)	Difference (Limits)	Flags	A or P
	M-131	VD-080310				
Perchlorate	60000 ug/L	60000 ug/L	0 (≤30)	-	-	-
Total dissolved solids	3200 mg/L	3200 mg/L	0 (≤30)	-	-	-

**2010 Annual Remedial Performance Sampling  
Wet Chemistry - Data Qualification Summary - SDG 339977**

SDG	Sample	Analyte	Flag	A or P	Reason
339977	I-Q I-C I-S I-L I-R I-B I-AR I-AB I-AA M-131 M-57A M-79 M-69 M-135 VD-080310	Perchlorate	J- (all detects) R (all non-detects)	P	Technical holding times
339977	EB080310V	Hexavalent chromium	J- (all detects) R (all non-detects)	P	Technical holding times

**2010 Annual Remedial Performance Sampling  
Wet Chemistry - Laboratory Blank Data Qualification Summary - SDG 339977**

No Sample Data Qualified in this SDG

**2010 Annual Remedial Performance Sampling  
Wet Chemistry - Field Blank Data Qualification Summary - SDG 339977**

No Sample Data Qualified in this SDG

Tronox Northgate Henderson

VALIDATION COMPLETENESS WORKSHEET

LDC #: 2420306  
 SDG #: 339977  
 Laboratory: MWH Laboratories

Stage 2 ~~B~~ A

9M4

Date: 12-15-10  
 Page: 1 of 1  
 Reviewer: MG  
 2nd Reviewer: LA

METHOD: (Analyte) Hexavalent Chromium (EPA SW846 Method 7196A), Perchlorate (EPA Method 314.0), TDS (EPA Method 160.1/SM2540C)

The samples listed below were reviewed for each of the following validation areas. Validation findings are noted in attached validation findings worksheets.

	Validation Area		Comments
I.	Technical holding times	SW	Sampling dates: 8-2-10 through 8-3-10
IIa.	Initial calibration	N	
IIb.	Calibration verification	N	
III.	Blanks	A	
IV	Matrix Spike/Matrix Spike Duplicates	A	MS/MSD (SDG: 339791)
V	Duplicates	A	DUP (SDG: 339791)
VI.	Laboratory control samples	A	LCS/LCSD
VII.	Sample result verification	N	
VIII.	Overall assessment of data	A	
IX.	Field duplicates	SW	D=10 +15
X	Field blanks	SW	EB = 16, FB = FB080210V (SDG: 339791)

Note: A = Acceptable  
 N = Not provided/applicable  
 SW = See worksheet  
 ND = No compounds detected  
 R = Rinsate  
 FB = Field blank  
 D = Duplicate  
 TB = Trip blank  
 EB = Equipment blank

Validated Samples: all water

1	I-Q	11	M-57A	21	EB080310VMS	31	
2	I-C	12	M-79	22	EB080310VMDS	32	
3	I-S	13	M-69	23		33	
4	I-L	14	M-135	24		34	
5	I-R	15	VD-080310	25		35	
6	I-B	16	EB080310V	26		36	
7	I-AR	17	I-CMS	27	PBW1	37	
8	I-AB	18	I-CMSD	28	PBW2	38	
9	I-AA	19	I-CDUP	29	PBW3	39	
10	M-131	20	M-79DUP	30	PBW4	40	

Notes: \_\_\_\_\_

**VALIDATION FINDINGS WORKSHEET  
Sample Specific Analysis Reference**

All circled methods are applicable to each sample.

Sample ID	Matrix	Parameter
1 → 15	W	pH (TDS) CI F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> (ClO <sub>4</sub> )
16	↓	pH (TDS) CI F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC (CR <sup>6+</sup> ) (ClO <sub>4</sub> )
QC 17, 18		pH TDS CI F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> (ClO <sub>4</sub> )
↓ 19, 20	↓	pH (TDS) CI F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
↓ 21, 22		pH TDS CI F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC (CR <sup>6+</sup> ) ClO <sub>4</sub>
		pH TDS CI F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
		pH TDS CI F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
		pH TDS CI F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
		pH TDS CI F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
		pH TDS CI F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
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		pH TDS CI F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
		pH TDS CI F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
		pH TDS CI F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
		pH TDS CI F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
		pH TDS CI F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>
	pH TDS CI F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>	
	pH TDS CI F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>	
	pH TDS CI F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>	
	pH TDS CI F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>	
	pH TDS CI F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>	
	pH TDS CI F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> ClO <sub>4</sub>	

Comments:

**VALIDATION FINDINGS WORKSHEET**  
**Technical Holding Times**

All circled dates have exceeded the technical holding time.

N N/A Were all samples preserved as applicable to each method?

N N/A Were all cooler temperatures within validation criteria?

Method:	314.0	7196 A					
Parameters:	C104	Cr VI					
Technical holding time:	28 days	24 hr					
Sample ID	Sampling date	Analysis date	Analysis date	Analysis date	Analysis date	Analysis date	Qualifier
1 → 8, 17, 18	8-2-10	10-13-10 (72 days)					J-/R/P
9	↓	10-14-10 (73 days)					
10	8-3-10	↓ (72 days)					
11 → 15	↓	10-13-10 (71 days)					
16	<sup>0208</sup> 8-3-10			<sup>1503</sup> 8-5-10 (61.00 hr)			
21	↓			↓			
22	↓			↓			

**VALIDATION FINDINGS WORKSHEET**  
**Field Blanks**

**METHOD:** Inorganics, EPA Method. See Cover  
 N/A Were field blanks identified in this SDG?  
 N/A Were target analytes detected in the field blanks?  
**Blank units:** ug/L. **Associated sample units:** ug/L  
**Sampling date:** 8/2/10. **Soil factor applied:** NA  
**Field blank type:** (circle one) Field Blank / Rinsate / Other.

Associated Samples: ClO4: 1-9 (>RL), Cr VI: none

Analyte	Blank ID	Action Limit	Sample Identification
	VD080210V	No Qual's.	
Cr VI (mg/L)	0.0060	0.06	
ClO4	1.63		

**Blank units:** ug/L. **Associated sample units:** ug/L  
**Sampling date:** 8/3/10. **Soil factor applied:** NA  
**Field blank type:** (circle one) Field Blank / Rinsate / Other EB

Associated Samples: 10-15 (>10x)

Analyte	Blank ID	Action Limit	Sample Identification
	16	No Qual's.	
ClO4	8.5	85	

CIRCLED RESULTS WERE NOT QUALIFIED. ALL RESULTS NOT CIRCLED WERE QUALIFIED BY THE FOLLOWING STATEMENT:  
 Samples with analyte concentrations within five times the associated field blank concentration are listed above, these sample results were qualified as not detected, "U".

LDC#: 2420306  
SDG#: See Cover

**VALIDATION FINDINGS WORKSHEET**  
**Field Duplicates**

Page: 1 of 1  
Reviewer: MG  
2nd Reviewer: [Signature]

Inorganics, Method See Cover

- N NA Were field duplicate pairs identified in this SDG?
- N NA Were target analytes detected in the field duplicate pairs?

Analyte	Concentration (mg/L)		RPD ( $\leq 30$ )	Difference	Limits	Qualification (Parent only)
	10	15				
Perchlorate (ug/L)	60000	60000	0			
TDS	3200	3200	0			

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## Laboratory Data Consultants, Inc. Data Validation Report

**Project/Site Name:** 2010 Annual Remedial Performance Sampling  
**Collection Date:** August 1 through August 2, 2010  
**LDC Report Date:** December 23, 2010  
**Matrix:** Water  
**Parameters:** Perchlorate  
**Validation Level:** Stage 2A  
**Laboratory:** MWH Laboratories, Inc.  
**Sample Delivery Group (SDG):** 339791

### Sample Identification

I-O	M-96
I-P	PC-54
I-H	PC-37
I-U	PC-71
I-T	PC-72
I-G	PC-73
I-F	M-23
I-N	VD080210
I-E	FB080210V
I-M	FB080210VMS
I-D	FB080210VMMSD
PC-123	
PC-124	
PC-125	
PC-126	
PC-127	
PC-128	
PC-129	
PC-130	
PC-132	

## Introduction

This data review covers 31 water samples listed on the cover sheet including dilutions and reanalysis as applicable. The analyses were per EPA Method 331.0 for Perchlorate.

This review follows a modified outline of the USEPA Contract Laboratory Program National Functional Guidelines for Inorganic Data Review (October 2004) and the EPA Region 9 Superfund Data Evaluation/Validation Guidance, NDEP guidance (May 2006).

A qualification summary table is provided at the end of this report if data has been qualified. Flags are classified as P (protocol) or A (advisory) to indicate whether the flag is due to a laboratory deviation from a specified protocol or is of technical advisory nature.

Raw data were not reviewed for this SDG. The review was based on QC data.

The following are definitions of the data qualifiers:

- J+ Data are qualified as estimated, with a high bias likely to occur. False positives or false negatives are unlikely to have been reported.
- J- Data are qualified as estimated, with a low bias likely to occur. False positives or false negatives are unlikely to have been reported.
- J Data are qualified as estimated; it is not possible to assess the direction of the potential bias. False positives or false negatives are unlikely to have been reported.
- U Indicates the compound or analyte was analyzed for but not detected at or above the stated limit.
- R Data are qualified as rejected. There is a significant potential for the reporting of false negatives or false positives.
- UJ Indicates the compound or analyte was analyzed for but not detected. The sample detection limit is an estimated value.
- A Indicates the finding is based upon technical validation criteria.
- P Indicates the finding is related to a protocol/contractual deviation.
- None Indicates the data was not significantly impacted by the finding, therefore qualification was not required.

## I. Technical Holding Times

All technical holding time requirements were met.

The chain-of-custodies were reviewed for documentation of cooler temperatures. All cooler temperatures met validation criteria.

## II. LC/MS Instrument Performance Check

Instrument performance check data were not reviewed for Stage 2A.

## III. Initial Calibration

Initial calibration data were not reviewed for Stage 2A.

## IV. Continuing Calibration

Calibration verification data were not reviewed for Stage 2A.

## V. Blanks

Method blanks were reviewed for each matrix as applicable. No perchlorate was found in the method blanks.

Sample FB080210V was identified as a field blank. No contaminant concentrations were found in this blank with the following exceptions:

Field Blank ID	Sampling Date	Analyte	Concentration	Associated Samples
FB080210V	8/2/10	Perchlorate	2 ug/L	I-O I-P I-H I-U I-T I-G I-F I-N I-E I-M I-D PC-54 PC-37 PC-71 PC-72 PC-73 M-23 VD080210

Sample concentrations were compared to concentrations detected in the field blanks as required by the QAPP. No sample data was qualified.

## **VI. Surrogate Spikes**

Surrogates were not required by the method.

## **VII. Matrix Spike/Matrix Spike Duplicates**

Matrix spike (MS) and matrix spike duplicate (MSD) samples were reviewed for each matrix as applicable. Percent recoveries (%R) and relative percent differences (RPD) were within QC limits.

## **VIII. Laboratory Control Samples (LCS)**

Laboratory control samples were reviewed for each matrix as applicable. Percent recoveries (%R) and relative percent differences (RPD) were within QC limits.

## **IX. Regional Quality Assurance and Quality Control**

Not applicable.

## **X. Internal Standards**

Internal standards data were not reviewed for Stage 2A.

## **XI. Target Compound Identifications**

Raw data were not reviewed for this SDG.

## **XII. Compound Quantitation and CRQLs**

Raw data were not reviewed for this SDG.

## **XIII. Tentatively Identified Compounds (TICs)**

Raw data were not reviewed for this SDG.

## **XIV. System Performance**

Raw data were not reviewed for this SDG.

## **XV. Overall Assessment**

Data flags are summarized at the end of this report if data has been qualified.

## XVI. Field Duplicates

Samples M-23 and VD080210 were identified as field duplicates. No perchlorate was detected in any of the samples with the following exceptions:

Analyte	Concentration (ug/L)		RPD (Limits)	Flags	A or P
	M-23	VD080210			
Perchlorate	460000	320000	36 ( $\leq 30$ )	J (all detects)	A

**2010 Annual Remedial Performance Sampling  
Perchlorate - Data Qualification Summary - SDG 339791**

SDG	Sample	Analyte	Flag	A or P	Reason
339791	M-23 VD080210	Perchlorate	J (all detects)	A	Field duplicates (RPD)

**2010 Annual Remedial Performance Sampling  
Perchlorate - Laboratory Blank Data Qualification Summary - SDG 339791**

No Sample Data Qualified in this SDG

**2010 Annual Remedial Performance Sampling  
Perchlorate - Field Blank Data Qualification Summary - SDG 339791**

No Sample Data Qualified in this SDG

LDC #: 24203N87

## VALIDATION COMPLETENESS WORKSHEET

Date: 12-15-10

SDG #: 339791

Stage 2B A

Page: 1 of 1

Laboratory: MWH Laboratories

Reviewer: MG

2nd Reviewer:         

METHOD: LC/MS Perchlorate (EPA Method 331.0)

The samples listed below were reviewed for each of the following validation areas. Validation findings are noted in attached validation findings worksheets.

	Validation Area		Comments
I.	Technical holding times	A	Sampling dates: 8-1-10 through 8-2-10
II.	GC/MS Instrument performance check	N	
III.	Initial calibration	N	
IV.	Continuing calibration/ICV	N	
V.	Blanks	A	
VI.	Surrogate spikes	N	
VII.	Matrix spike/Matrix spike duplicates	A	MS/MSD
VIII.	Laboratory control samples	A	LCS/LCSD
IX.	Regional Quality Assurance and Quality Control	N	
X.	Internal standards	N	
XI.	Target compound identification	N	
XII.	Compound quantitation/CRQLs	N	
XIII.	Tentatively identified compounds (TICs)	N	
XIV.	System performance	N	
XV.	Overall assessment of data	A	
XVI.	Field duplicates	SW	D = 27 + 28
XVII.	Field blanks	SW	FB = 29

Note: A = Acceptable  
N = Not provided/applicable  
SW = See worksheet

ND = No compounds detected  
R = Rinsate  
FB = Field blank

D = Duplicate  
TB = Trip blank  
EB = Equipment blank

Validated Samples:

all water

1	I-O	11	I-D	21	M-96	31	FB080210VMSD
2	I-P	12	PC-123	22	PC-54	32	PBW1
3	I-H	13	PC-124	23	PC-37	33	PBW2
4	I-U	14	PC-125	24	PC-71	34	
5	I-T	15	PC-126	25	PC-72	35	
6	I-G	16	PC-127	26	PC-73	36	
7	I-F	17	PC-128	27	M-23	37	
8	I-N	18	PC-129	28	VD080210	38	
9	I-E	19	PC-130	29	FB080210V	39	
10	I-M	20	PC-132	30	FB080210VMS	40	

# VALIDATION FINDINGS WORKSHEET

## Field Blanks

**METHOD:** Inorganics, EPA Method See Cover  
 N N/A Were field blanks identified in this SDG?  
 N N/A Were target analytes detected in the field blanks?  
**Blank units:** ug/L **Associated sample units:** ug/L  
**Sampling date:** 8/2/10 **Soil factor applied:** NA  
**Field blank type:** (circle one)  Field Blank  Rinsate / Other:

Associated Samples: ClO4: 1-11, 22-28 (>10x)

Analyte	Blank ID	Action Limit	No Qual's	Sample Identification
	29		No Qual's	
ClO4	2	20		

CIRCLED RESULTS WERE NOT QUALIFIED. ALL RESULTS NOT CIRCLED WERE QUALIFIED BY THE FOLLOWING STATEMENT:  
 Samples with analyte concentrations within five times the associated field blank concentration are listed above, these sample results were qualified as not detected, "U".



LDC#: 24203N87

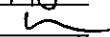
SDG#: See Cover

### VALIDATION FINDINGS WORKSHEET

#### Field Duplicates

Page: 1 of 1

Reviewer: MG

2nd Reviewer: 

Inorganics, Method See Cover

Y  N  NA

Were field duplicate pairs identified in this SDG?

Y  N  NA

Were target analytes detected in the field duplicate pairs?

Analyte	Concentration (ug/L)		RPD ( $\leq 30$ )	Difference	Limits	Qualification (Parent only)
	27	28				
Perchlorate	460000	320000	36			J dets/ A fd

V:\FIELD DUPLICATES\FD\_inorganic\24203N87.wpd

**Laboratory Data Consultants, Inc.  
Data Validation Report**

**Project/Site Name:** 2010 Annual Remedial Performance Sampling  
**Collection Date:** August 2 through August 3, 2010  
**LDC Report Date:** December 23, 2010  
**Matrix:** Water  
**Parameters:** Perchlorate  
**Validation Level:** Stage 2A  
**Laboratory:** MWH Laboratories, Inc.  
**Sample Delivery Group (SDG):** 339977

**Sample Identification**

I-Q  
I-C  
I-S  
I-L  
I-R  
I-B  
I-AR  
I-AB  
I-AA  
M-131  
M-57A  
M-79  
M-69  
M-135  
VD-080310  
EB080310V  
EB080310VMS  
EB080310VMSD

## Introduction

This data review covers 18 water samples listed on the cover sheet including dilutions and reanalysis as applicable. The analyses were per EPA Method 331.0 for Perchlorate.

This review follows a modified outline of the USEPA Contract Laboratory Program National Functional Guidelines for Inorganic Data Review (October 2004) and the EPA Region 9 Superfund Data Evaluation/Validation Guidance, NDEP guidance (May 2006).

A qualification summary table is provided at the end of this report if data has been qualified. Flags are classified as P (protocol) or A (advisory) to indicate whether the flag is due to a laboratory deviation from a specified protocol or is of technical advisory nature.

Raw data were not reviewed for this SDG. The review was based on QC data.

The following are definitions of the data qualifiers:

**J+** Data are qualified as estimated, with a high bias likely to occur. False positives or false negatives are unlikely to have been reported.

**J-** Data are qualified as estimated, with a low bias likely to occur. False positives or false negatives are unlikely to have been reported.

**J** Data are qualified as estimated; it is not possible to assess the direction of the potential bias. False positives or false negatives are unlikely to have been reported.

**U** Indicates the compound or analyte was analyzed for but not detected at or above the stated limit.

**R** Data are qualified as rejected. There is a significant potential for the reporting of false negatives or false positives.

**UJ** Indicates the compound or analyte was analyzed for but not detected. The sample detection limit is an estimated value.

**A** Indicates the finding is based upon technical validation criteria.

**P** Indicates the finding is related to a protocol/contractual deviation.

**None** Indicates the data was not significantly impacted by the finding, therefore qualification was not required.

## I. Technical Holding Times

All technical holding time requirements were met.

The chain-of-custodies were reviewed for documentation of cooler temperatures. All cooler temperatures met validation criteria.

## II. LC/MS Instrument Performance Check

Instrument performance check data were not reviewed for Stage 2A.

## III. Initial Calibration

Initial calibration data were not reviewed for Stage 2A.

## IV. Continuing Calibration

Calibration verification data were not reviewed for Stage 2A.

## V. Blanks

Method blanks were reviewed for each matrix as applicable. No perchlorate was found in the method blanks.

Sample EB080310V was identified as an equipment blank. No contaminant concentrations were found in this blank with the following exceptions:

Equipment Blank ID	Sampling Date	Analyte	Concentration	Associated Samples
EB080310V	8/3/10	Perchlorate	10 ug/L	M-131 M-57A M-79 M-69 M-135 VD-080310

Sample FB080210V (from SDG 339791) was identified as a field blank. No contaminant concentrations were found in this blank with the following exceptions:

Field Blank ID	Sampling Date	Analyte	Concentration	Associated Samples
FB080210V	8/2/10	Perchlorate	2 ug/L	I-Q I-C I-S I-L I-R I-B I-AR I-AB I-AA

Sample concentrations were compared to concentrations detected in the field blanks as required by the QAPP. No sample data was qualified.

#### **VI. Surrogate Spikes**

Surrogates were not required by the method.

#### **VII. Matrix Spike/Matrix Spike Duplicates**

Matrix spike (MS) and matrix spike duplicate (MSD) samples were reviewed for each matrix as applicable. Percent recoveries (%R) and relative percent differences (RPD) were within QC limits.

#### **VIII. Laboratory Control Samples (LCS)**

Laboratory control samples were reviewed for each matrix as applicable. Percent recoveries (%R) and relative percent differences (RPD) were within QC limits.

#### **IX. Regional Quality Assurance and Quality Control**

Not applicable.

#### **X. Internal Standards**

Internal standards data were not reviewed for Stage 2A.

#### **XI. Target Compound Identifications**

Raw data were not reviewed for this SDG.

#### **XII. Compound Quantitation and CRQLs**

Raw data were not reviewed for this SDG.

#### **XIII. Tentatively Identified Compounds (TICs)**

Raw data were not reviewed for this SDG.

#### **XIV. System Performance**

Raw data were not reviewed for this SDG.

#### **XV. Overall Assessment**

Data flags are summarized at the end of this report if data has been qualified.

## XVI. Field Duplicates

Samples M-131 and VD-080310 were identified as field duplicates. No perchlorate was detected in any of the samples with the following exceptions:

Compound	Concentration (ug/L)		RPD (Limits)	Flags	A or P
	M-131	VD-080310			
Perchlorate	64000	66000	3 ( $\leq 30$ )	-	-

**2010 Annual Remedial Performance Sampling  
Perchlorate - Data Qualification Summary - SDG 339977**

No Sample Data Qualified in this SDG

**2010 Annual Remedial Performance Sampling  
Perchlorate - Laboratory Blank Data Qualification Summary - SDG 339977**

No Sample Data Qualified in this SDG

**2010 Annual Remedial Performance Sampling  
Perchlorate - Field Blank Data Qualification Summary - SDG 339977**

No Sample Data Qualified in this SDG

LDC #: 24203087

## VALIDATION COMPLETENESS WORKSHEET

Date: 10-15-10

SDG #: 339977

Stage 2B A

Page: 1 of 1

Laboratory: MWH Laboratories

Reviewer: MG

2nd Reviewer: ✓

METHOD: LC/MS Perchlorate (EPA Method 331.0)

The samples listed below were reviewed for each of the following validation areas. Validation findings are noted in attached validation findings worksheets.

	Validation Area		Comments
I.	Technical holding times	A	Sampling dates: 8-2-10 through 8-3-10
II.	GC/MS Instrument performance check	N	
III.	Initial calibration	N	
IV.	Continuing calibration/ICV	N	
V.	Blanks	A	
VI.	Surrogate spikes	N	
VII.	Matrix spike/Matrix spike duplicates	A	MS/MSD (SDG: 339791)
VIII.	Laboratory control samples	A	LCS/LCSD
IX.	Regional Quality Assurance and Quality Control	N	
X.	Internal standards	N	
XI.	Target compound identification	N	
XII.	Compound quantitation/CRQLs	N	
XIII.	Tentatively identified compounds (TICs)	N	
XIV.	System performance	N	
XV.	Overall assessment of data	A	
XVI.	Field duplicates	SW	D = 10 + 15
XVII.	Field blanks	SW	EB = 16 FB = FB080210V (SDG: 339791)

Note: A = Acceptable  
N = Not provided/applicable  
SW = See worksheet

ND = No compounds detected  
R = Rinsate  
FB = Field blank

D = Duplicate  
TB = Trip blank  
EB = Equipment blank

Validated Samples:  
all water

1	I-Q	11	M-57A	21	31
2	I-C	12	M-79	22	32
3	I-S	13	M-69	23	33
4	I-L	14	M-135	24	34
5	I-R	15	VD-080310	25	35
6	I-B	16	EB080310V	26	36
7	I-AR	17	EB080310VMS	27	37
8	I-AB	18	EB080310VMSD	28	38
9	I-AA	19	PBW1	29	39
10	M-131	20	PBW2	30	40



**VALIDATION FINDINGS WORKSHEET**  
**Field Blanks**

**METHOD:** Inorganics, EPA Method See Cover  
 N/A Were field blanks identified in this SDG?  
 N/A Were target analytes detected in the field blanks?  
**Blank units:** ug/L **Associated sample units:** ug/L  
**Sampling date:** 8/2/10 **Soil factor applied:** NA  
**Field blank type:** (circle one) Field Blank / Rinsate / Other: \_\_\_\_\_

Associated Samples: CIO4: 1-9 (>10x)

Analyte	Blank ID	Action Limit	Sample Identification
	VD080210V	No Qual's.	
CIO4	2	20	

**Blank units:** ug/L **Associated sample units:** ug/L  
**Sampling date:** 8/3/10 **Soil factor applied:** NA  
**Field blank type:** (circle one) Field Blank / Rinsate / Other EB

Associated Samples: 10-15 (>10x)

Analyte	Blank ID	Action Limit	Sample Identification
	16	No Qual's.	
CIO4	10	100	

CIRCLED RESULTS WERE NOT QUALIFIED. ALL RESULTS NOT CIRCLED WERE QUALIFIED BY THE FOLLOWING STATEMENT:  
 Samples with analyte concentrations within five times the associated field blank concentration are listed above, these sample results were qualified as not detected, "U".

LDC#: 24203087  
SDG#: See Cover

**VALIDATION FINDINGS WORKSHEET**  
**Field Duplicates**

Page: 1 of 1  
Reviewer: MG  
2nd Reviewer: L

Inorganics, Method See Cover

- Y N NA Were field duplicate pairs identified in this SDG?  
 Y N NA Were target analytes detected in the field duplicate pairs?

Analyte	Concentration (ug/L)		RPD ( $\leq 30$ )	Difference	Limits	Qualification (Parent only)
	10	15				
Perchlorate	64000	66000	3			

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