

**Tronox LLC Facility, 2008 Phase B Investigation, Henderson, Nevada
Data Validation Reports
LDC# 21257**

Dioxins/Dibenzofurans

LDC

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

Project/Site Name: Tronox LLC Facility, 2008 Phase B Investigation,
Henderson, Nevada

Collection Date: June 30, 2008

LDC Report Date: August 21, 2009

Matrix: Soil

Parameters: Dioxins/Dibenzofurans

Validation Level: Stage 2B

Laboratory: Columbia Analytical Services, Inc.

Sample Delivery Group (SDG): E0800626

Sample Identification

SA207-0.5B

SA207-0.5BDL

Introduction

This data review covers 2 soil samples listed on the cover sheet including dilutions and reanalysis as applicable. The analyses were per EPA SW 846 Method 8290 for Polychlorinated Dioxins/Dibenzofurans.

This review follows the Standard Operating Procedures (SOP) 40, Data Review/Validation (BRC 2009), the Quality Assurance Project Plan Tronox LLC Facility, Henderson, Nevada (June 2009), NDEP guidance (May 2006), and USEPA Contract Laboratory Program National Functional Guidelines for Polychlorinated Dioxins/Dibenzofurans Data Review (September 2005) as there are no current guidelines for the method 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.

Blank results are summarized in Section V.

Field duplicates are summarized in Section XIV.

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.
- B The analytical result may be a false positive totally attributable to blank contamination. This qualifier is applicable to radiochemistry analysis only.
- JB The analytical result may be biased high and partially attributable to blank contamination. This qualifier is applicable to radiochemistry analysis only.
- JK The analytical result is an estimated maximum possible concentration (EMPC).
- X The analytical result is not used for reporting because a more accurate and precise result is reported in its place.
- J-TDS The analytical result is estimated based on failure of the Total Dissolved Solids (TDS) correctness check performed in accordance with the Standard Method 1030E.
- J-CAB The analytical result is estimated based on failure of the cation-anion balance correctness check performed in accordance with Standard Method 1030E.
- J-TDS & CAB The analytical result is unreliable based on the failure of the cation-anion balance and TDS correctness check performed in accordance with standard Method 1030E.
- 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. HRGC/HRMS Instrument Performance Check

Instrument performance was checked at the required daily frequency.

Retention time windows were established for all homologues. The chromatographic resolution between 2,3,7,8-TCDD and peaks representing any other unlabeled TCDD isomer was less than or equal to 25%.

III. Initial Calibration

A five point initial calibration was performed as required by the method.

Percent relative standard deviations (%RSD) were less than or equal to 20.0% for unlabeled compounds and less than or equal to 30.0% for labeled compounds.

The ion abundance ratios for all PCDDs and PCDFs were within validation criteria.

IV. Routine Calibration (Continuing)

Routine calibration was performed at the required frequencies.

All of the routine calibration percent differences (%D) between the initial calibration RRF and the routine calibration RRF were less than or equal to 20.0% for unlabeled compounds and less than or equal to 30.0% for labeled compounds.

The ion abundance ratios for all PCDDs and PCDFs were within validation criteria.

V. Blanks

Method blanks were reviewed for each matrix as applicable. No polychlorinated dioxin/dibenzofuran contaminants were found in the method blanks with the following exceptions:

| Method Blank ID | Extraction Date | Compound | Concentration | Associated Samples |
|-----------------|-----------------|--|---|--------------------------------|
| EQ0800286-01 | 7/3/08 | 1,2,3,4,6,7,8-HpCDD OCDD 1,2,3,4,6,7,8-HpCDF Total HpCDD Total HpCDF | 0.135 ng/Kg 0.487 ng/Kg 0.0867 ng/Kg 0.351 ng/Kg 0.0867 ng/Kg | All samples in SDG E0800626 |

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

No field blanks were identified in this SDG.

VI. Matrix Spike/Matrix Spike Duplicates

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. 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 with the following exceptions:

| LCS ID (Associated Samples) | Compound | LCS %R (Limits) | LCSD %R (Limits) | RPD (Limits) | Flag | A or P |
|---|--------------|--------------------|---------------------|-----------------|------------------|--------|
| EQ0800286-02LCS/D (All samples in SDG E0800626) | 2,3,7,8-TCDD | 137 (87-135) | 136 (87-135) | - | J+ (all detects) | P |

VIII. Regional Quality Assurance and Quality Control

Not applicable.

IX. Internal Standards

All internal standard recoveries were within QC limits.

X. Target Compound Identifications

Raw data were not reviewed for this SDG.

XI. Project Quantitation Limit

All project quantitation limits were within validation criteria with the following exceptions:

| Sample | Compound | Finding | Criteria | Flag | A or P |
|--------------|--|---|---|--|--------|
| SA207-0.5B | 2,3,7,8-TCDD 1,2,3,7,8-PeCDD 1,2,3,4,7,8-HxCDD 1,2,3,6,7,8-HxCDD 1,2,3,7,8,9-HxCDD 1,2,3,4,6,7,8-HpCDD OCDD 2,3,7,8-TCDF 2,3,4,7,8-PeCDF 1,2,3,4,7,8-HxCDF 1,2,3,6,7,8-HxCDF 2,3,4,6,7,8-HxCDF 1,2,3,7,8,9-HxCDF 1,2,3,4,6,7,8-HpCDF 1,2,3,4,7,8,9-HpCDF OCDF | Sample result exceeded calibration range. | Reported result should be within calibration range. | J (all detects) J (all detects) J (all detects) J (all detects) J (all detects) J (all detects) J (all detects) J (all detects) J (all detects) J (all detects) J (all detects) J (all detects) J (all detects) J (all detects) J (all detects) J (all detects) | A |
| SA207-0.5BDL | 2,3,7,8-TCDF 1,2,3,4,6,7,8-HpCDF OCDF | Sample result exceeded calibration range. | Reported result should be within calibration range. | J (all detects) J (all detects) J (all detects) | A |

All compounds reported below the PQL were qualified as follows:

| Sample | Finding | Flag | A or P |
|-----------------------------|---------------------------------------|-----------------|--------|
| All samples in SDG E0800626 | All compounds reported below the PQL. | J (all detects) | A |

All compounds reported as EMPC were qualified as follows:

| Sample | Finding | Flag | A or P |
|-----------------------------|--|------------------|--------|
| All samples in SDG E0800626 | All compounds reported as estimated maximum possible concentration (EMPC). | JK (all detects) | A |

Raw data were not reviewed for this SDG.

XII. System Performance

Raw data were not reviewed for this SDG.

XIII. Overall Assessment of Data

The overall assessment of data was acceptable. In the case where more than one result was reported for an individual sample, the least technically acceptable results were rejected as follows:

| Sample | Compound | Flag | A or P |
|--------------|------------------------|------|--------|
| SA207-0.5B | All TCL compounds | X | A |
| SA207-0.5BDL | 2,3,7,8-TCDF from DB-5 | X | A |

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.

**Tronox LLC Facility, 2008 Phase B Investigation, Henderson, Nevada
Dioxins/Dibenzofurans - Data Qualification Summary - SDG E0800626**

| SDG | Sample | Compound | Flag | A or P | Reason (Code) |
|----------|----------------------------|--|---|--------|-------------------------------------|
| E0800626 | SA207-0.5B SA207-0.5BDL | 2,3,7,8-TCDD | J+ (all detects) | P | Laboratory control samples (%R) (l) |
| E0800626 | SA207-0.5B | 2,3,7,8-TCDD 1,2,3,7,8-PeCDD 1,2,3,4,7,8-HxCDD 1,2,3,6,7,8-HxCDD 1,2,3,7,8,9-HxCDD 1,2,3,4,6,7,8-HpCDD OCDD 2,3,7,8-TCDF 2,3,4,7,8-PeCDF 1,2,3,4,7,8-HxCDF 1,2,3,6,7,8-HxCDF 2,3,4,6,7,8-HxCDF 1,2,3,7,8,9-HxCDF 1,2,3,4,6,7,8-HpCDF 1,2,3,4,7,8,9-HpCDF OCDF | J (all detects) J (all detects) J (all detects) J (all detects) J (all detects) J (all detects) J (all detects) J (all detects) J (all detects) J (all detects) J (all detects) J (all detects) J (all detects) J (all detects) J (all detects) J (all detects) J (all detects) J (all detects) J (all detects) | A | Project Quantitation Limit (e) |
| E0800626 | SA207-0.5BDL | 2,3,7,8-TCDF 1,2,3,4,6,7,8-HpCDF OCDF | J (all detects) J (all detects) J (all detects) | A | Project Quantitation Limit (e) |
| E0800626 | SA207-0.5B SA207-0.5BDL | All compounds reported below the PQL. | J (all detects) | A | Project Quantitation Limit (sp) |
| E0800626 | SA207-0.5B SA207-0.5BDL | All compounds reported as EMPC | JK (all detects) | A | Project Quantitation Limit (k) |
| E0800626 | SA207-0.5B | All TCL compounds | X | A | Overall assessment of data (o) |
| E0800626 | SA207-0.5BDL | 2,3,7,8-TCDF from DB-5 | X | A | Overall assessment of data (o) |

**Tronox LLC Facility, 2008 Phase B Investigation, Henderson, Nevada
Dioxins/Dibenzofurans - Laboratory Blank Data Qualification Summary - SDG E0800626**

No Sample Data Qualified in this SDG

**Tronox LLC Facility, 2008 Phase B Investigation, Henderson, Nevada
Dioxins/Dibenzofurans - Field Blank Data Qualification Summary - SDG E0800626**

No Sample Data Qualified in this SDG

Tronox Northgate Henderson

VALIDATION COMPLETENESS WORKSHEET

Stage 2B

LDC #: 21257N21
 SDG #: E0800626
 Laboratory: Columbia Analytical Services

Date: 8/14/09
 Page: 1 of 1
 Reviewer: [Signature]
 2nd Reviewer: [Signature]

METHOD: HRGC/HRMS Dioxins/Dibenzofurans (EPA SW 846 Method 8290)

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: 6/30/08 |
| II. | HRGC/HRMS Instrument performance check | A | |
| III. | Initial calibration | A | |
| IV. | Routine calibration/MSV | A | |
| V. | Blanks | SW | |
| VI. | Matrix spike/Matrix spike duplicates | N | |
| VII. | Laboratory control samples | SW | LCs/D |
| VIII. | Regional quality assurance and quality control | N | |
| IX. | Internal standards | A | |
| X. | Target compound identifications | N | |
| XI. | Compound quantitation and CRQLs | SW | |
| XII. | System performance | N | |
| XIII. | Overall assessment of data | SW | |
| XIV. | Field duplicates | N | |
| XV. | 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:

| | | | | | | | |
|----|------------|---|-----|--------------|-----|---------|----|
| 1 | SA207-0.5B | S | 11' | EA0800286-01 | 21 | U216TAT | 31 |
| 2 | ↓ DL | S | 12 | | 22' | 1 768 | 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: _____

VALIDATION FINDINGS WORKSHEET

METHOD: HRGC/HRMS Dioxins/Dibenzofurans (EPA SW 846 Method 8290)

| | | | | |
|----------------------|------------------------|------------------------|------------------------|----------------|
| A. 2,3,7,8-TCDD | F. 1,2,3,4,6,7,8-HpCDD | K. 1,2,3,4,7,8-HxCDF | P. 1,2,3,4,7,8,9-HpCDF | U. Total HpCDD |
| B. 1,2,3,7,8-PeCDD | G. OCDD | L. 1,2,3,6,7,8-HxCDF | Q. OCDF | V. Total TCDF |
| C. 1,2,3,4,7,8-HxCDD | H. 2,3,7,8-TCDF | M. 2,3,4,6,7,8-HxCDF | R. Total TCDD | W. Total PeCDF |
| D. 1,2,3,6,7,8-HxCDD | I. 1,2,3,7,8-PeCDF | N. 1,2,3,7,8,9-HxCDF | S. Total PeCDD | X. Total HxCDF |
| E. 1,2,3,7,8,9-HxCDD | J. 2,3,4,7,8-PeCDF | O. 1,2,3,4,6,7,8-HpCDF | T. Total HxCDD | Y. Total HpCDF |

Notes:

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

Project/Site Name: Tronox LLC Facility, 2008 Phase B Investigation,
Henderson, Nevada

Collection Date: July 2, 2008

LDC Report Date: August 24, 2009

Matrix: Soil

Parameters: Dioxins/Dibenzofurans

Validation Level: Stage 2B

Laboratory: Columbia Analytical Services, Inc.

Sample Delivery Group (SDG): E0800632

Sample Identification

SA181-0.5B

Introduction

This data review covers one soil sample listed on the cover sheet including dilutions and reanalysis as applicable. The analyses were per EPA SW 846 Method 8290 for Polychlorinated Dioxins/Dibenzofurans.

This review follows the Standard Operating Procedures (SOP) 40, Data Review/Validation (BRC 2009), the Quality Assurance Project Plan Tronox LLC Facility, Henderson, Nevada (June 2009), NDEP guidance (May 2006), and USEPA Contract Laboratory Program National Functional Guidelines for Polychlorinated Dioxins/Dibenzofurans Data Review (September 2005) as there are no current guidelines for the method 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.

Blank results are summarized in Section V.

Field duplicates are summarized in Section XIV.

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.
- B The analytical result may be a false positive totally attributable to blank contamination. This qualifier is applicable to radiochemistry analysis only.
- JB The analytical result may be biased high and partially attributable to blank contamination. This qualifier is applicable to radiochemistry analysis only.
- JK The analytical result is an estimated maximum possible concentration (EMPC).
- X The analytical result is not used for reporting because a more accurate and precise result is reported in its place.
- J-TDS The analytical result is estimated based on failure of the Total Dissolved Solids (TDS) correctness check performed in accordance with the Standard Method 1030E.
- J-CAB The analytical result is estimated based on failure of the cation-anion balance correctness check performed in accordance with Standard Method 1030E.
- J-TDS & CAB The analytical result is unreliable based on the failure of the cation-anion balance and TDS correctness check performed in accordance with standard Method 1030E.
- 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. HRGC/HRMS Instrument Performance Check

Instrument performance was checked at the required daily frequency.

Retention time windows were established for all homologues. The chromatographic resolution between 2,3,7,8-TCDD and peaks representing any other unlabeled TCDD isomer was less than or equal to 25%.

III. Initial Calibration

A five point initial calibration was performed as required by the method.

Percent relative standard deviations (%RSD) were less than or equal to 20.0% for unlabeled compounds and less than or equal to 30.0% for labeled compounds.

The ion abundance ratios for all PCDDs and PCDFs were within validation criteria.

IV. Routine Calibration (Continuing)

Routine calibration was performed at the required frequencies.

All of the routine calibration percent differences (%D) between the initial calibration RRF and the routine calibration RRF were less than or equal to 20.0% for unlabeled compounds and less than or equal to 30.0% for labeled compounds.

The ion abundance ratios for all PCDDs and PCDFs were within validation criteria.

V. Blanks

Method blanks were reviewed for each matrix as applicable. No polychlorinated dioxin/dibenzofuran contaminants were found in the method blanks with the following exceptions:

| Method Blank ID | Extraction Date | Compound | Concentration | Associated Samples |
|-----------------|-----------------|--|---|--------------------------------|
| EQ0800286-01 | 7/34/08 | 1,2,3,4,6,7,8-HpCDD OCDD 1,2,3,4,6,7,8-HpCDF Total HpCDD Total HpCDF | 0.135 ng/Kg 0.487 ng/Kg 0.0867 ng/Kg 0.351 ng/Kg 0.0867 ng/Kg | All samples in SDG E0800632 |

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 | Compound | Reported Concentration | Modified Final Concentration |
|------------|--|--|---|
| SA181-0.5B | 1,2,3,4,6,7,8-HpCDD OCDD Total HpCDD | 0.387 ng/Kg 2.28 ng/Kg 0.799 ng/Kg | 0.387U ng/Kg 2.28U ng/Kg 0.799U ng/Kg |

No field blanks were identified in this SDG.

VI. Matrix Spike/Matrix Spike Duplicates

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. 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 with the following exceptions:

| LCS ID (Associated Samples) | Compound | LCS %R (Limits) | LCSD %R (Limits) | RPD (Limits) | Flag | A or P |
|---|--------------|--------------------|---------------------|-----------------|------------------|--------|
| EQ0800286-02LCS/D (All samples in SDG E0800632) | 2,3,7,8-TCDD | 137 (87-135) | 136 (87-135) | - | J+ (all detects) | P |

VIII. Regional Quality Assurance and Quality Control

Not applicable.

IX. Internal Standards

All internal standard recoveries were within QC limits.

X. Target Compound Identifications

Raw data were not reviewed for this SDG.

XI. Project Quantitation Limit

All compounds reported below the PQL were qualified as follows:

| Sample | Finding | Flag | A or P |
|-----------------------------|---------------------------------------|-----------------|--------|
| All samples in SDG E0800632 | All compounds reported below the PQL. | J (all detects) | A |

All compounds reported as EMPC were qualified as follows:

| Sample | Finding | Flag | A or P |
|-----------------------------|--|------------------|--------|
| All samples in SDG E0800632 | All compounds reported as estimated maximum possible concentration (EMPC). | JK (all detects) | A |

Raw data were not reviewed for this SDG.

XII. System Performance

Raw data were not reviewed for this SDG.

XIII. Overall Assessment of Data

The overall assessment of data was acceptable. In the case where more than one result was reported for an individual sample, the least technically acceptable results were rejected as follows:

| Sample | Compound | Flag | A or P |
|------------|------------------------|------|--------|
| SA181-0.5B | 2,3,7,8-TCDF from DB-5 | X | A |

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.

**Tronox LLC Facility, 2008 Phase B Investigation, Henderson, Nevada
Dioxins/Dibenzofurans - Data Qualification Summary - SDG E0800632**

| SDG | Sample | Compound | Flag | A or P | Reason (Code) |
|----------|------------|--------------------------------------|------------------|--------|-------------------------------------|
| E0800632 | SA181-0.5B | 2,3,7,8-TCDD | J+ (all detects) | P | Laboratory control samples (%R) (l) |
| E0800632 | SA181-0.5B | All compounds reported below the PQL | J (all detects) | A | Project Quantitation Limit (sp) |
| E0800632 | SA181-0.5B | All compounds reported as EMPC | JK (all detects) | A | Project Quantitation Limit (k) |
| E0800632 | SA181-0.5B | 2,3,7,8-TCDF from DB-5 | X | A | Overall assessment of data (o) |

**Tronox LLC Facility, 2008 Phase B Investigation, Henderson, Nevada
Dioxins/Dibenzofurans - Laboratory Blank Data Qualification Summary - SDG E0800632**

| SDG | Sample | Compound | Modified Final Concentration | A or P | Code |
|----------|------------|--|---|--------|------|
| E0800632 | SA181-0.5B | 1,2,3,4,6,7,8-HpCDD OCDD Total HpCDD | 0.387U ng/Kg 2.28U ng/Kg 0.799U ng/Kg | A | bl |

**Tronox LLC Facility, 2008 Phase B Investigation, Henderson, Nevada
Dioxins/Dibenzofurans - Field Blank Data Qualification Summary - SDG E0800632**

No Sample Data Qualified in this SDG

Tronox Northgate Henderson

VALIDATION COMPLETENESS WORKSHEET

Stage 2B

LDC #: 21257021
 SDG #: E0800632
 Laboratory: Columbia Analytical Services

Date: 8/14/09
 Page: 1 of 1
 Reviewer: [Signature]
 2nd Reviewer: [Signature]

METHOD: HRGC/HRMS Dioxins/Dibenzofurans (EPA SW 846 Method 8290)

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: 7/2/09 |
| II. | HRGC/HRMS Instrument performance check | A | |
| III. | Initial calibration | A | |
| IV. | Routine calibration/IDV | A | |
| V. | Blanks | SW | |
| VI. | Matrix spike/Matrix spike duplicates | N | direct purified |
| VII. | Laboratory control samples | SW | CS/D |
| VIII. | Regional quality assurance and quality control | N | |
| IX. | Internal standards | A | |
| X. | Target compound identifications | N | |
| XI. | Compound quantitation and CRQLs | SW | All ZMPC results - JK |
| XII. | System performance | N | |
| XIII. | Overall assessment of data | SW | |
| 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:

| | | | | | | | |
|----|------------|---|----|---------------|----|---------|----|
| 1 | SA181-0.5B | S | 11 | ZR0800-286-01 | 21 | U216T9T | 31 |
| 2 | | | 12 | | 22 | U216T68 | 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: _____

VALIDATION FINDINGS WORKSHEET

METHOD: HRGC/HRMS Dioxins/Dibenzofurans (EPA SW 846 Method 8290)

| | | | | |
|----------------------|------------------------|------------------------|------------------------|----------------|
| A. 2,3,7,8-TCDD | F. 1,2,3,4,6,7,8-HpCDD | K. 1,2,3,4,7,8-HxCDF | P. 1,2,3,4,7,8,9-HpCDF | U. Total HpCDD |
| B. 1,2,3,7,8-PeCDD | G. OCDD | L. 1,2,3,6,7,8-HxCDF | Q. OCDF | V. Total TCDF |
| C. 1,2,3,4,7,8-HxCDD | H. 2,3,7,8-TCDF | M. 2,3,4,6,7,8-HxCDF | R. Total TCDD | W. Total PeCDF |
| D. 1,2,3,6,7,8-HxCDD | I. 1,2,3,7,8-PeCDF | N. 1,2,3,7,8,9-HxCDF | S. Total PeCDD | X. Total HxCDF |
| E. 1,2,3,7,8,9-HxCDD | J. 2,3,4,7,8-PeCDF | O. 1,2,3,4,6,7,8-HpCDF | T. Total HxCDD | Y. Total HpCDF |

Notes:

VALIDATION FINDINGS WORKSHEET
Blanks

METHOD: HRGC/HRMS Dioxins/Dibenzofurans (EPA Method 8290)

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

- N N/A Were all samples associated with a method blank?
- Y N N/A Was a method blank analyzed for each matrix?
- Y N N/A Was the blank contaminated? If yes, please see qualification below.

Blank extraction date: 7/3/08 Blank analysis date: 7/13/08

Conc. units: ng/kg Associated Samples: M (6)

| Compound | Blank ID | Sample Identification |
|----------|--------------|-----------------------|
| | EA0502286-01 | 1 |
| F | 0.135 | 0.387/u |
| F | 0.487 | 2.28/u |
| O | 0.0867 | 0.09 |
| U | 0.351 | 0.799/u |
| Y | 0.0267 | 0.90 |
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Blank extraction date: _____ Blank analysis date: _____
 Conc. units: _____ Associated Samples: _____

| Compound | Blank ID | Sample Identification |
|----------|----------|-----------------------|
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CIRCLED RESULTS WERE NOT QUALIFIED. ALL RESULTS NOT CIRCLED WERE QUALIFIED BY THE FOLLOWING STATEMENT:
 All contaminants within five times the method blank concentration were qualified as not detected, "U".

VALIDATION FINDINGS WORKSHEET Overall Assessment of Data

METHOD: HRGC/HRMS Dioxins/Dibenzofurans (EPA SW 846 Method 8290)

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

All available information pertaining to the data were reviewed using professional judgement to compliment the determination of the overall quality of the data.

(Y) N N/A Was the overall quality and usability of the data acceptable?

| # | Date | Sample ID | Finding | Associated Samples | Qualifications |
|---|------|-----------|------------------|--------------------|----------------|
| | | <i>1</i> | <i>1 on DB-5</i> | <i>1</i> | <i>X (0)</i> |
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Comments: _____

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

Project/Site Name: Tronox LLC Facility, 2008 Phase B Investigation,
Henderson, Nevada

Collection Date: July 7 through July 8, 2008

LDC Report Date: August 26, 2009

Matrix: Soil

Parameters: Dioxins/Dibenzofurans

Validation Level: Stage 2B

Laboratory: Columbia Analytical Services, Inc.

Sample Delivery Group (SDG): E0800639

Sample Identification

SA47-0.5B
SA183-0.5B

All samples in this SDG were analyzed for screening purposes only.

Introduction

This data review covers 2 soil samples listed on the cover sheet including dilutions and reanalysis as applicable. The analyses were per EPA SW 846 Method 8290 for Polychlorinated Dioxins/Dibenzofurans.

This review follows the Standard Operating Procedures (SOP) 40, Data Review/Validation (BRC 2009), the Quality Assurance Project Plan Tronox LLC Facility, Henderson, Nevada (June 2009), NDEP guidance (May 2006), and USEPA Contract Laboratory Program National Functional Guidelines for Polychlorinated Dioxins/Dibenzofurans Data Review (September 2005) as there are no current guidelines for the method 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.

Blank results are summarized in Section V.

Field duplicates are summarized in Section XIV.

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.
- B The analytical result may be a false positive totally attributable to blank contamination. This qualifier is applicable to radiochemistry analysis only.
- JB The analytical result may be biased high and partially attributable to blank contamination. This qualifier is applicable to radiochemistry analysis only.
- JK The analytical result is an estimated maximum possible concentration (EMPC).
- X The analytical result is not used for reporting because a more accurate and precise result is reported in its place.
- J-TDS The analytical result is estimated based on failure of the Total Dissolved Solids (TDS) correctness check performed in accordance with the Standard Method 1030E.
- J-CAB The analytical result is estimated based on failure of the cation-anion balance correctness check performed in accordance with Standard Method 1030E.
- J-TDS & CAB The analytical result is unreliable based on the failure of the cation-anion balance and TDS correctness check performed in accordance with standard Method 1030E.
- 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. HRGC/HRMS Instrument Performance Check

Instrument performance check data were not reviewed for this SDG.

III. Initial Calibration

Initial calibration data were not reviewed for this SDG.

IV. Routine Calibration (Continuing)

Routine calibration data were not reviewed for this SDG.

V. Blanks

Method blanks were reviewed for each matrix as applicable. No polychlorinated dioxin/dibenzofuran contaminants were found in the method blanks with the following exceptions:

| Method Blank ID | Extraction Date | Compound | Concentration | Associated Samples |
|-----------------|-----------------|---|---|--------------------------------|
| EQ0800301-01 | 7/17/08 | 1,2,3,4,6,7,8-HpCDD OCDD 1,2,3,4,7,8-HxCDF 1,2,3,6,7,8-HxCDF 1,2,3,4,6,7,8-HpCDF 1,2,3,4,7,8,9-HpCDF OCDF | 0.354 ng/Kg 1.47 ng/Kg 0.364 ng/Kg 0.202 ng/Kg 0.825 ng/Kg 0.354 ng/Kg 2.42 ng/Kg | All samples in SDG E0800639 |

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

No field blanks were identified in this SDG.

VI. Matrix Spike/Matrix Spike Duplicates

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. 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 with the following exceptions:

| LCS ID (Associated Samples) | Compound | LCS %R (Limits) | LCSD %R (Limits) | RPD (Limits) | Flag | A or P |
|---|--------------|--------------------|---------------------|-----------------|------------------|--------|
| EQ0800301-02LCS/D (All samples in SDG E0800639) | 2,3,7,8-TCDD | 137 (87-135) | 137 (87-135) | - | J+ (all detects) | P |

VIII. Regional Quality Assurance and Quality Control

Not applicable.

IX. Internal Standards

Internal standards data were not reviewed for this SDG.

X. Target Compound Identifications

Raw data were not reviewed for this SDG.

XI. Project Quantitation Limit

All compounds reported below the PQL were qualified as follows:

| Sample | Finding | Flag | A or P |
|-----------------------------|---------------------------------------|-----------------|--------|
| All samples in SDG E0800639 | All compounds reported below the PQL. | J (all detects) | A |

Raw data were not reviewed for this SDG.

XII. System Performance

Raw data were not reviewed for this SDG.

XIII. Overall Assessment of Data

Samples SA47-0.5B and SA183-0.5B were analyzed for screening purposes only.

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.

**Tronox LLC Facility, 2008 Phase B Investigation, Henderson, Nevada
Dioxins/Dibenzofurans - Data Qualification Summary - SDG E0800639**

| SDG | Sample | Compound | Flag | A or P | Reason (Code) |
|----------|-------------------------|---------------------------------------|------------------|--------|-------------------------------------|
| E0800639 | SA47-0.5B SA183-0.5B | 2,3,7,8-TCDD | J+ (all detects) | P | Laboratory control samples (%R) (I) |
| E0800639 | SA47-0.5B SA183-0.5B | All compounds reported below the PQL. | J (all detects) | A | Project Quantitation Limit (sp) |

**Tronox LLC Facility, 2008 Phase B Investigation, Henderson, Nevada
Dioxins/Dibenzofurans - Laboratory Blank Data Qualification Summary - SDG E0800639**

No Sample Data Qualified in this SDG

**Tronox LLC Facility, 2008 Phase B Investigation, Henderson, Nevada
Dioxins/Dibenzofurans - Field Blank Data Qualification Summary - SDG E0800639**

No Sample Data Qualified in this SDG

LDC #: 21257P21
 SDG #: E0800639
 Laboratory: Columbia Analytical Services

Tronox Northgate Henderson
VALIDATION COMPLETENESS WORKSHEET
 Stage 2B

Date: 8/14/09
 Page: 1 of 1
 Reviewer: [Signature]
 2nd Reviewer: [Signature]

METHOD: HRGC/HRMS Dioxins/Dibenzofurans (EPA SW 846 Method 8290)

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>7/7-8/08</u> |
| II. | HRGC/HRMS Instrument performance check | N | |
| III. | Initial calibration | N | |
| IV. | Routine calibration/ICV | N | |
| V. | Blanks | SW | |
| VI. | Matrix spike/Matrix spike duplicates | N | |
| VII. | Laboratory control samples | SW | <u>LCS/D</u> |
| VIII. | Regional quality assurance and quality control | N | |
| IX. | Internal standards | N | |
| X. | Target compound identifications | N | |
| XI. | Compound quantitation and CRQLs | SW | |
| XII. | System performance | N | |
| XIII. | Overall assessment of data | A | |
| XIV. | Field duplicates | N | |
| XV. | 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:

| | | | | | | | | | |
|----|------------|---|---|----|---------------------|----|--|----|--|
| 1 | SA47-0.5B | * | S | 11 | <u>ZQ0800301-01</u> | 21 | | 31 | |
| 2 | SA183-0.5B | * | ✓ | 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: * Samples were analyzed for screening only

VALIDATION FINDINGS WORKSHEET
Overall Assessment of Data

Page: 1 of 1
 Reviewer: _____
 2nd Reviewer: _____

LDC #: 225731
 SDG #: 225731

METHOD: HRGC/HRMS Dioxins/Dibenzofurans (EPA SW 846 Method 8290)

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

All available information pertaining to the data were reviewed using professional judgement to compliment the determination of the overall quality of the data.

N N/A Was the overall quality and usability of the data acceptable?

| # | Date | Sample ID | Finding | Associated Samples | Qualifications |
|---|------|-----------|--|--------------------|----------------|
| | | MM | Samples were screening analyzed for screening only. data was not validated . | MM | Text |
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Comments: _____

VALIDATION FINDINGS WORKSHEET

METHOD: HRGC/HRMS Dioxins/Dibenzofurans (EPA SW 846 Method 8290)

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|----------------------|------------------------|------------------------|------------------------|----------------|
| A. 2,3,7,8-TCDD | F. 1,2,3,4,6,7,8-HpCDD | K. 1,2,3,4,7,8-HxCDF | P. 1,2,3,4,7,8,9-HpCDF | U. Total HpCDD |
| B. 1,2,3,7,8-PeCDD | G. OCDD | L. 1,2,3,6,7,8-HxCDF | Q. OCDF | V. Total TCDF |
| C. 1,2,3,4,7,8-HxCDD | H. 2,3,7,8-TCDF | M. 2,3,4,6,7,8-HxCDF | R. Total TCDD | W. Total PeCDF |
| D. 1,2,3,6,7,8-HxCDD | I. 1,2,3,7,8-PeCDF | N. 1,2,3,7,8,9-HxCDF | S. Total PeCDD | X. Total HxCDF |
| E. 1,2,3,7,8,9-HxCDD | J. 2,3,4,7,8-PeCDF | O. 1,2,3,4,6,7,8-HpCDF | T. Total HxCDD | Y. Total HpCDF |

Notes:

VALIDATION FINDINGS WORKSHEET
Blanks

METHOD: HRGC/HRMS Dioxins/Dibenzofurans (EPA Method 8290)

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

N N/A Were all samples associated with a method blank?

N N/A Was a method blank analyzed for each matrix?

N N/A Was the blank contaminated? If yes, please see qualification below.

Blank extraction date: 7/17/08 Blank analysis date: 7/28/08

Conc. units: ng/kg Associated Samples: nd (>5x)

| Compound | Blank ID | Sample Identification |
|----------|-------------------------|-----------------------|
| | EA0800301-01 | |
| F | 0.354 | |
| G | 1.47 | |
| K | 0.364 | |
| L | 0.202 | |
| O | 0.825 | |
| P | 0.354 | |
| Q | 2.42 | |

Blank extraction date: _____ Blank analysis date: _____

Conc. units: _____ Associated Samples: _____

| Compound | Blank ID | Sample Identification |
|--------------|------------------|-----------------------|
| U | 0.503 | |
| X | 0.364 | |
| Y | 0.469 | |
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CIRCLED RESULTS WERE NOT QUALIFIED. ALL RESULTS NOT CIRCLED WERE QUALIFIED BY THE FOLLOWING STATEMENT:
 All contaminants within five times the method blank concentration were qualified as not detected, "U".

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

Project/Site Name: Tronox LLC Facility, 2008 Phase B Investigation,
Henderson, Nevada

Collection Date: July 7 through July 8, 2008

LDC Report Date: August 21, 2009

Matrix: Soil

Parameters: Dioxins/Dibenzofurans

Validation Level: Stage 2B

Laboratory: Columbia Analytical Services, Inc.

Sample Delivery Group (SDG): E0800640

Sample Identification

SA67-0.5B

RSAN2-0.5B

Introduction

This data review covers 2 soil samples listed on the cover sheet including dilutions and reanalysis as applicable. The analyses were per EPA SW 846 Method 8290 for Polychlorinated Dioxins/Dibenzofurans.

This review follows the Standard Operating Procedures (SOP) 40, Data Review/Validation (BRC 2009), the Quality Assurance Project Plan Tronox LLC Facility, Henderson, Nevada (June 2009), NDEP guidance (May 2006), and USEPA Contract Laboratory Program National Functional Guidelines for Polychlorinated Dioxins/Dibenzofurans Data Review (September 2005) as there are no current guidelines for the method 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.

Blank results are summarized in Section V.

Field duplicates are summarized in Section XIV.

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.
- B The analytical result may be a false positive totally attributable to blank contamination. This qualifier is applicable to radiochemistry analysis only.
- JB The analytical result may be biased high and partially attributable to blank contamination. This qualifier is applicable to radiochemistry analysis only.
- JK The analytical result is an estimated maximum possible concentration (EMPC).
- X The analytical result is not used for reporting because a more accurate and precise result is reported in its place.
- J-TDS The analytical result is estimated based on failure of the Total Dissolved Solids (TDS) correctness check performed in accordance with the Standard Method 1030E.
- J-CAB The analytical result is estimated based on failure of the cation-anion balance correctness check performed in accordance with Standard Method 1030E.
- J-TDS & CAB The analytical result is unreliable based on the failure of the cation-anion balance and TDS correctness check performed in accordance with standard Method 1030E.
- 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. HRGC/HRMS Instrument Performance Check

Instrument performance was checked at the required daily frequency.

Retention time windows were established for all homologues. The chromatographic resolution between 2,3,7,8-TCDD and peaks representing any other unlabeled TCDD isomer was less than or equal to 25%.

III. Initial Calibration

A five point initial calibration was performed as required by the method.

Percent relative standard deviations (%RSD) were less than or equal to 20.0% for unlabeled compounds and less than or equal to 30.0% for labeled compounds.

The ion abundance ratios for all PCDDs and PCDFs were within validation criteria.

IV. Routine Calibration (Continuing)

Routine calibration was performed at the required frequencies.

All of the routine calibration percent differences (%D) between the initial calibration RRF and the routine calibration RRF were less than or equal to 20.0% for unlabeled compounds and less than or equal to 30.0% for labeled compounds.

The ion abundance ratios for all PCDDs and PCDFs were within validation criteria.

V. Blanks

Method blanks were reviewed for each matrix as applicable. No polychlorinated dioxin/dibenzofuran contaminants were found in the method blanks with the following exceptions:

| Method Blank ID | Extraction Date | Compound | Concentration | Associated Samples |
|-----------------|-----------------|----------|---------------|--------------------|
| EQ0800312-01 | 7/23/08 | OCDD | 0.850 ng/Kg | SA67-0.5B |

| Method Blank ID | Extraction Date | Compound | Concentration | Associated Samples |
|-----------------|-----------------|--|--|--------------------|
| EQ0800301-01 | 7/17/08 | 1,2,3,4,6,7,8-HpCDD OCDD 1,2,3,4,7,8-HxCDF 1,2,3,6,7,8-HxCDF 1,2,3,4,6,7,8-HpCDF 1,2,3,4,7,8,9-HpCDF OCDF Total HpCDD Total HxCDF Total HpCDF | 0.354 ng/Kg 1.47 ng/Kg 0.364 ng/Kg 0.202 ng/Kg 0.825 ng/Kg 0.354 ng/Kg 2.42 ng/Kg 0.503 ng/Kg 0.364 ng/Kg 0.469 ng/Kg | RSAN2-0.5B |

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

No field blanks were identified in this SDG.

VI. Matrix Spike/Matrix Spike Duplicates

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. 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 with the following exceptions:

| LCS ID (Associated Samples) | Compound | LCS %R (Limits) | LCSD %R (Limits) | RPD (Limits) | Flag | A or P |
|---|--------------|--------------------|---------------------|-----------------|------------------|--------|
| EQ0800301-02LCS/D (RSAN2-0.5B EQ0800301-01) | 2,3,7,8-TCDD | 137 (87-135) | 137 (87-135) | - | J+ (all detects) | P |
| EQ0800312-02LCS/D (SA67-0.5B EQ0800312-01) | 2,3,7,8-TCDD | 136 (87-135) | 139 (87-135) | - | J+ (all detects) | P |

VIII. Regional Quality Assurance and Quality Control

Not applicable.

IX. Internal Standards

All internal standard recoveries were within QC limits with the following exceptions:

| Sample | Internal Standards | %R (Limits) | Compound | Flag | A or P |
|--------------|---|----------------------------|---|--|--------|
| RSAN2-0.5B | ¹³ C-OCDD ¹³ C-1,2,3,4,6,7,8-HpCDF | 36 (40-135) 39 (40-135) | OCDD 1,2,3,4,6,7,8-HpCDF 1,2,3,4,7,8,9-HpCDF OCDF Total HpCDF | J (all detects) UJ (all non-detects) | P |
| EQ0800312-01 | ¹³ C-OCDD | 25 (40-135) | OCDD OCDF | J (all detects) UJ (all non-detects) J (all detects) UJ (all non-detects) | P |

X. Target Compound Identifications

Raw data were not reviewed for this SDG.

XI. Project Quantitation Limit

All project quantitation limits were within validation criteria with the following exceptions:

| Sample | Compound | Finding | Criteria | Flag | A or P |
|------------|---|---|---|---|--------|
| RSAN2-0.5B | 1,2,3,4,6,7,8-HpCDD OCDD 2,3,7,8-TCDF 1,2,3,4,7,8-HxCDF 1,2,3,4,6,7,8-HpCDF | Sample result exceeded calibration range. | Reported result should be within calibration range. | J (all detects) J (all detects) J (all detects) J (all detects) J (all detects) | P |

All compounds reported below the PQL were qualified as follows:

| Sample | Finding | Flag | A or P |
|-----------------------------|---------------------------------------|-----------------|--------|
| All samples in SDG E0800640 | All compounds reported below the PQL. | J (all detects) | A |

All compounds reported as EMPC were qualified as follows:

| Sample | Finding | Flag | A or P |
|-----------------------------|--|------------------|--------|
| All samples in SDG E0800640 | All compounds reported as estimated maximum possible concentration (EMPC). | JK (all detects) | A |

Raw data were not reviewed for this SDG.

XII. System Performance

Raw data were not reviewed for this SDG.

XIII. Overall Assessment of Data

The overall assessment of data was acceptable. In the case where more than one result was reported for an individual sample, the least technically acceptable results were rejected as follows:

| Sample | Compound | Flag | A or P |
|-------------------------|------------------------|------|--------|
| SA67-0.5B RSAN2-0.5B | 2,3,7,8-TCDF from DB-5 | X | A |

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.

**Tronox LLC Facility, 2008 Phase B Investigation, Henderson, Nevada
Dioxins/Dibenzofurans - Data Qualification Summary - SDG E0800640**

| SDG | Sample | Compound | Flag | A or P | Reason (Code) |
|----------|-------------------------|---|---|--------|-------------------------------------|
| E0800640 | SA67-0.5B RSAN2-0.5B | 2,3,7,8-TCDD | J+ (all detects) | P | Laboratory control samples (%R) (l) |
| E0800640 | RSAN2-0.5B | OCDD 1,2,3,4,6,7,8-HpCDF 1,2,3,4,7,8,9-HpCDF OCDF Total HpCDF | J (all detects) J (all detects) J (all detects) J (all detects) J (all detects) | A | Internal standards (area) (i) |
| E0800640 | RSAN2-0.5B | 1,2,3,4,6,7,8-HpCDD OCDD 2,3,7,8-TCDF 1,2,3,4,7,8-HxCDF 1,2,3,4,6,7,8-HpCDF | J (all detects) J (all detects) J (all detects) J (all detects) J (all detects) | P | Project Quantitation Limit (e) |
| E0800640 | SA67-0.5B RSAN2-0.5B | All compounds reported below the PQL. | J (all detects) | A | Project Quantitation Limit (sp) |
| E0800640 | SA67-0.5B RSAN2-0.5B | All compounds reported as EMPC | JK (all detects) | A | Project Quantitation Limit (k) |
| E0800640 | SA67-0.5B RSAN2-0.5B | 2,3,7,8-TCDF from DB-5 | X | A | Overall assessment of data (o) |

**Tronox LLC Facility, 2008 Phase B Investigation, Henderson, Nevada
Dioxins/Dibenzofurans - Laboratory Blank Data Qualification Summary - SDG E0800640**

No Sample Data Qualified in this SDG

**Tronox LLC Facility, 2008 Phase B Investigation, Henderson, Nevada
Dioxins/Dibenzofurans - Field Blank Data Qualification Summary - SDG E0800640**

No Sample Data Qualified in this SDG

Tronox Northgate Henderson

VALIDATION COMPLETENESS WORKSHEET

Stage 2B

LDC #: 21257Q21
 SDG #: E0800640
 Laboratory: Columbia Analytical Services

Date: 8/14/09
 Page: 1 of 1
 Reviewer: [Signature]
 2nd Reviewer: [Signature]

METHOD: HRGC/HRMS Dioxins/Dibenzofurans (EPA SW 846 Method 8290)

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: 7/7 - 8/08 |
| II. | HRGC/HRMS Instrument performance check | A |
| III. | Initial calibration | A |
| IV. | Routine calibration/ICV | A |
| V. | Blanks | W |
| VI. | Matrix spike/Matrix spike duplicates | N direct spiked |
| VII. | Laboratory control samples | W CS/O |
| VIII. | Regional quality assurance and quality control | N |
| IX. | Internal standards | W |
| X. | Target compound identifications | N |
| XI. | Compound quantitation and CRQLs | SW |
| XII. | System performance | N |
| XIII. | Overall assessment of data | W |
| XIV. | Field duplicates | N |
| XV. | 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:

| | | | | | | | | |
|----|------------|---|----|---|--------------|----|------------|----|
| 1 | SA67-0.5B | S | 11 | 2 | 200800301-01 | 21 | P100148 | 31 |
| 2 | RSAN2-0.5B | ↓ | 12 | 1 | 200800312-01 | 22 | C15238 # 2 | 32 |
| 3 | | | 13 | | | 23 | M216901 | 33 |
| 4 | | | 14 | | | 24 | C15225 # 2 | 34 |
| 5 | | | 15 | | | 25 | P100128 | 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

METHOD: HRGC/HRMS Dioxins/Dibenzofurans (EPA SW 846 Method 8290)

| | | | | |
|----------------------|------------------------|------------------------|------------------------|----------------|
| A. 2,3,7,8-TCDD | F. 1,2,3,4,6,7,8-HpCDD | K. 1,2,3,4,7,8-HxCDF | P. 1,2,3,4,7,8,9-HpCDF | U. Total HpCDD |
| B. 1,2,3,7,8-PeCDD | G. OCDD | L. 1,2,3,6,7,8-HxCDF | Q. OCDF | V. Total TCDF |
| C. 1,2,3,4,7,8-HxCDD | H. 2,3,7,8-TCDF | M. 2,3,4,6,7,8-HxCDF | R. Total TCDD | W. Total PeCDF |
| D. 1,2,3,6,7,8-HxCDD | I. 1,2,3,7,8-PeCDF | N. 1,2,3,7,8,9-HxCDF | S. Total PeCDD | X. Total HxCDF |
| E. 1,2,3,7,8,9-HxCDD | J. 2,3,4,7,8-PeCDF | O. 1,2,3,4,6,7,8-HpCDF | T. Total HxCDD | Y. Total HpCDF |

Notes:

VALIDATION FINDINGS WORKSHEET
Blanks

METHOD: HRGC/HRMS Dioxins/Dibenzofurans (EPA Method 8290)

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

- Y N N/A Were all samples associated with a method blank?
- Y N N/A Was a method blank analyzed for each matrix?
- Y N N/A Was the blank contaminated? If yes, please see qualification below.

Blank extraction date: 7/17/08 **Blank analysis date:** 7/22/08

Conc. units: ng/kg **Associated Samples:** 2 (25 X)

| Compound | Blank ID | Sample Identification |
|----------|--------------|-----------------------|
| | EA1800301-01 | |
| F | 0.354 | |
| G | 1.47 | |
| K | 0.364 | |
| L | 0.202 | |
| O | 0.825 | |
| P | 0.354 | |
| Q | 2.42 | |

Blank extraction date: _____ **Blank analysis date:** _____ **Associated Samples:** _____

| Compound | Blank ID | Sample Identification |
|----------|----------|-----------------------|
| | | |
| U | 0.503 | |
| X | 0.364 | |
| X | 0.469 | |
| | | |
| | | |
| | | |
| | | |
| | | |
| | | |

CIRCLED RESULTS WERE NOT QUALIFIED. ALL RESULTS NOT CIRCLED WERE QUALIFIED BY THE FOLLOWING STATEMENT:
 All contaminants within five times the method blank concentration were qualified as not detected, "U".

VALIDATION FINDINGS WORKSHEET
Internal Standards

METHOD: HRGC/HRMS Dioxins/Dibenzofurans (EPA SW 846 Method 8290)

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

Y/N N/A Are all internal standard recoveries were within the 40-135% criteria?

Y/N N/A Was the S/N ratio all internal standard peaks > 10?

| # | Date | Lab ID/Reference | Internal Standard | % Recovery (Limit: 40-135%) | Qualifications |
|---|------|--------------------|-------------------|-----------------------------|---------------------------------------|
| | | <u>2</u> | <u>I</u> | <u>36</u> (40-135) | <u>✓</u> <u>MP (i)</u> |
| | | | <u>9</u> | <u>39</u> () | <u>✓</u> |
| | | | | () | |
| | | | | () | |
| | | | | () | |
| | | | | () | |
| | | <u>88080312-01</u> | <u>I</u> | <u>25</u> (40-135) | <u>✓</u> <u>MP (i)</u> (<u>9.8</u>) |
| | | | | () | |
| | | | | () | |
| | | | | () | |
| | | | | () | |
| | | | | () | |
| | | | | () | |
| | | | | () | |
| | | | | () | |
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| | | | | () | |
| | | | | () | |
| | | | | () | |

| | Internal Standards | Check Standard Used | Internal Standards | Check Standard Used |
|----|-------------------------------------|---------------------|--------------------|-----------------------------------|
| A. | ¹³ C-2,3,7,8-TCDF | | I. | ¹³ C-OCDD |
| B. | ¹³ C-2,3,7,8-TCDD | | K. | ¹³ C-1,2,3,4-TCDD |
| C. | ¹³ C-1,2,3,7,8-PeCDF | | L. | ¹³ C-1,2,3,7,8,9-HxCDD |
| D. | ¹³ C-1,2,3,7,8-PeCDD | | M. | |
| E. | ¹³ C-1,2,3,4,7,8-HxCDF | | N. | |
| F. | ¹³ C-1,2,3,6,7,8-HxCDD | | O. | |
| G. | ¹³ C-1,2,3,4,6,7,8-HpCDF | | P. | |
| H. | ¹³ C-1,2,3,4,6,7,8-HpCDD | | | |

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

Project/Site Name: Tronox LLC Facility, 2008 Phase B Investigation,
Henderson, Nevada

Collection Date: July 10 through July 11, 2008

LDC Report Date: September 10, 2009

Matrix: Soil

Parameters: Dioxins/Dibenzofurans

Validation Level: Stage 2B

Laboratory: Columbia Analytical Services, Inc.

Sample Delivery Group (SDG): E0800661

Sample Identification

RSAJ8-0.5B
RSAI7-0.5B
RSAJ8-0.5BDL
RSAI7-0.5BDL

Introduction

This data review covers 4 soil samples listed on the cover sheet including dilutions and reanalysis as applicable. The analyses were per EPA SW 846 Method 8290 for Polychlorinated Dioxins/Dibenzofurans.

This review follows the Standard Operating Procedures (SOP) 40, Data Review/Validation (BRC 2009), the Quality Assurance Project Plan Tronox LLC Facility, Henderson, Nevada (June 2009), NDEP guidance (May 2006), and USEPA Contract Laboratory Program National Functional Guidelines for Polychlorinated Dioxins/Dibenzofurans Data Review (September 2005) as there are no current guidelines for the method 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.

Blank results are summarized in Section V.

Field duplicates are summarized in Section XIV.

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.
- B The analytical result may be a false positive totally attributable to blank contamination. This qualifier is applicable to radiochemistry analysis only.
- JB The analytical result may be biased high and partially attributable to blank contamination. This qualifier is applicable to radiochemistry analysis only.
- JK The analytical result is an estimated maximum possible concentration (EMPC).
- X The analytical result is not used for reporting because a more accurate and precise result is reported in its place.
- J-TDS The analytical result is estimated based on failure of the Total Dissolved Solids (TDS) correctness check performed in accordance with the Standard Method 1030E.
- J-CAB The analytical result is estimated based on failure of the cation-anion balance correctness check performed in accordance with Standard Method 1030E.
- J-TDS & CAB The analytical result is unreliable based on the failure of the cation-anion balance and TDS correctness check performed in accordance with standard Method 1030E.
- 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. HRGC/HRMS Instrument Performance Check

Instrument performance was checked at the required daily frequency.

Retention time windows were established for all homologues. The chromatographic resolution between 2,3,7,8-TCDD and peaks representing any other unlabeled TCDD isomer was less than or equal to 25%.

III. Initial Calibration

A five point initial calibration was performed as required by the method.

Percent relative standard deviations (%RSD) were less than or equal to 20.0% for unlabeled compounds and less than or equal to 30.0% for labeled compounds.

The ion abundance ratios for all PCDDs and PCDFs were within validation criteria.

IV. Routine Calibration (Continuing)

Routine calibration was performed at the required frequencies.

All of the routine calibration percent differences (%D) between the initial calibration RRF and the routine calibration RRF were less than or equal to 20.0% for unlabeled compounds and less than or equal to 30.0% for labeled compounds with the following exceptions:

| Date | Compound | %D | Associated Samples | Affected Compound | Flag | A or P |
|---------|--|-------------------------|--------------------|--|--|--------|
| 7/23/08 | ¹³ C-1,2,3,7,8-PeCDD ¹³ C-OCDD ¹³ C-1,2,3,7,8-PeCDF | 43.69 49.67 39.28 | EQ0800299-01 | 1,2,3,7,8-PeCDD OCDD 1,2,3,7,8-PeCDF | J- (all detects) UJ (all non-detects) | P |

The ion abundance ratios for all PCDDs and PCDFs were within validation criteria.

V. Blanks

Method blanks were reviewed for each matrix as applicable. No polychlorinated dioxin/dibenzofuran contaminants were found in the method blanks with the following exceptions:

| Method Blank ID | Extraction Date | Compound | Concentration | Associated Samples |
|-----------------|-----------------|---|--|----------------------------|
| EQ0800294-01 | 7/14/08 | 1,2,3,4,6,7,8-HpCDD OCDD 2,3,7,8-TCDF 1,2,3,7,8-PeCDF 1,2,3,4,7,8-HxCDF 1,2,3,6,7,8-HxCDF 1,2,3,4,6,7,8-HpCDF 1,2,3,4,7,8,9-HpCDF OCDF Total PeCDD Total HpCDD Total TCDF Total PeCDF Total HxCDF Total HpCDF | 0.265 ng/Kg 1.10 ng/Kg 0.581 ng/Kg 0.276 ng/Kg 0.340 ng/Kg 0.229 ng/Kg 0.405 ng/Kg 0.177 ng/Kg 0.796 ng/Kg 0.0845 ng/Kg 0.547 ng/Kg 2.13 ng/Kg 1.67 ng/Kg 1.14 ng/Kg 0.603 ng/Kg | RSAI7-0.5B RSAI7-0.5BDL |
| EQ0800299-01 | 7/16/08 | 1,2,3,4,6,7,8-HpCDD OCDD 1,2,3,4,6,7,8-HpCDF OCDF Total HpCDD Total HpCDF | 1.1004 ng/Kg 6.9141 ng/Kg 1.750 ng/Kg 4.050 ng/Kg 1.349 ng/Kg 1.750 ng/Kg | RSAJ8-0.5B |

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

No field blanks were identified in this SDG.

VI. Matrix Spike/Matrix Spike Duplicates

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. Laboratory Control Samples (LCS)

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

| LCS ID | Compound | %R (Limits) | Associated Samples | Flag | A or P |
|-----------------|-------------------------------------|------------------------------|----------------------------|--------------------------------------|--------|
| EQ0800299-03LCS | 2,3,7,8-TCDD 1,2,3,4,6,7,8-HpCDF | 144 (87-135) 144 (91-131) | RSAJ8-0.5B EQ0800299-01 | J+ (all detects) J+ (all detects) | P |

VIII. Regional Quality Assurance and Quality Control

Not applicable.

IX. Internal Standards

All internal standard recoveries were within QC limits with the following exceptions:

| Sample | Internal Standards | %R (Limits) | Compound | Flag | A or P |
|--------------|--|---|--|--|--------|
| RSAJ8-0.5B | ¹³ C-OCDD | 38 (40-135) | OCDD OCDF | J (all detects) UJ (all non-detects) J (all detects) UJ (all non-detects) | P |
| RSAI7-0.5B | ¹³ C-2,3,7,8-TCDF ¹³ C-1,2,3,4,7,8-HxCDF ¹³ C-1,2,3,6,7,8-HxCDD ¹³ C-1,2,3,4,6,7,8-HpCDF ¹³ C-1,2,3,4,6,7,8-HpCDD ¹³ C-OCDD | 12 (40-135) 25 (40-135) 27 (40-135) 15 (40-135) 20 (40-135) 12 (40-135) | 2,3,7,8-TCDF 1,2,3,4,7,8-HxCDD 1,2,3,6,7,8-HxCDD 1,2,3,7,8,9-HxCDD 1,2,3,4,6,7,8-HpCDD OCDD 1,2,3,4,7,8-HxCDF 1,2,3,6,7,8-HxCDF 2,3,4,6,7,8-HxCDF 1,2,3,7,8,9-HxCDF 1,2,3,4,6,7,8-HpCDF 1,2,3,4,7,8,9-HpCDF OCDF Total HxCDD Total HpCDD Total TCDF Total HxCDF Total HpCDF | J (all detects) UJ (all non-detects) | P |
| RSAI7-0.5BDL | ¹³ C-2,3,7,8-TCDF ¹³ C-1,2,3,4,7,8-HxCDF ¹³ C-1,2,3,6,7,8-HxCDD ¹³ C-1,2,3,4,6,7,8-HpCDF ¹³ C-1,2,3,4,6,7,8-HpCDD ¹³ C-OCDD | 16 (40-135) 24 (40-135) 30 (40-135) 17 (40-135) 23 (40-135) 11 (40-135) | 2,3,7,8-TCDF 1,2,3,4,7,8-HxCDD 1,2,3,6,7,8-HxCDD 1,2,3,7,8,9-HxCDD 1,2,3,4,6,7,8-HpCDD OCDD 1,2,3,4,7,8-HxCDF 1,2,3,6,7,8-HxCDF 2,3,4,6,7,8-HxCDF 1,2,3,7,8,9-HxCDF 1,2,3,4,6,7,8-HpCDF 1,2,3,4,7,8,9-HpCDF OCDF Total HxCDD Total HpCDD Total TCDF Total HxCDF Total HpCDF | J (all detects) UJ (all non-detects) | P |
| EQ0800299-01 | ¹³ C-2,3,7,8-TCDF ¹³ C-2,3,7,8-TCDD ¹³ C-1,2,3,7,8-PeCDF ¹³ C-1,2,3,7,8-PeCDD ¹³ C-1,2,3,4,7,8-HxCDF ¹³ C-1,2,3,6,7,8-HxCDD ¹³ C-1,2,3,4,6,7,8-HpCDF ¹³ C-1,2,3,4,6,7,8-HpCDD ¹³ C-OCDD | 10.61 (40-135) 11.53 (40-135) 12.27 (40-135) 12.28 (40-135) 10.12 (40-135) 11.00 (40-135) 10.20 (40-135) 11.23 (40-135) 9.32 (40-135) | All TCL compounds | J (all detects) UJ (all non-detects) | P |

X. Target Compound Identifications

Raw data were not reviewed for this SDG.

XI. Project Quantitation Limit

All project quantitation limits were within validation criteria with the following exceptions:

| Sample | Compound | Finding | Criteria | Flag | A or P |
|--------------|--|---|---|--|--------|
| RSAJ8-0.5B | 2,3,7,8-TCDF 1,2,3,4,7,8-HxCDF 1,2,3,4,6,7,8-HpCDF OCDF | Sample result exceeded calibration range. | Reported result should be within calibration range. | J (all detects) J (all detects) J (all detects) J (all detects) | A |
| RSAI7-0.5B | 2,3,7,8-TCDD 1,2,3,7,8-PeCDD 1,2,3,4,7,8-HxCDD 1,2,3,6,7,8-HxCDD 1,2,3,7,8,9-HxCDD 1,2,3,4,6,7,8-HpCDD OCDD 2,3,7,8-TCDF 2,3,4,7,8-PeCDF 1,2,3,4,7,8-HxCDF 1,2,3,6,7,8-HxCDF 2,3,4,6,7,8-HxCDF 1,2,3,7,8,9-HxCDF 1,2,3,4,6,7,8-HpCDF 1,2,3,4,7,8,9-HpCDF OCDF | Sample result exceeded calibration range. | Reported result should be within calibration range. | J (all detects) | A |
| RSAI7-0.5BDL | 2,3,7,8-TCDF 1,2,3,4,7,8-HxCDF 1,2,3,4,6,7,8-HpCDF 1,2,3,4,7,8,9-HpCDF OCDF | Sample result exceeded calibration range. | Reported result should be within calibration range. | J (all detects) J (all detects) J (all detects) J (all detects) J (all detects) | A |

All compounds reported below the PQL were qualified as follows:

| Sample | Finding | Flag | A or P |
|-----------------------------|---------------------------------------|-----------------|--------|
| All samples in SDG E0800661 | All compounds reported below the PQL. | J (all detects) | A |

All compounds reported as EMPC were qualified as follows:

| Sample | Finding | Flag | A or P |
|-----------------------------|--|------------------|--------|
| All samples in SDG E0800661 | All compounds reported as estimated maximum possible concentration (EMPC). | JK (all detects) | A |

Raw data were not reviewed for this SDG.

XII. System Performance

Raw data were not reviewed for this SDG.

XIII. Overall Assessment of Data

The overall assessment of data was acceptable. In the case where more than one result was reported for an individual sample, the least technically acceptable results were rejected as follows:

| Sample | Compound | Flag | A or P |
|---------------|--|-------------|---------------|
| RSAJ8-0.5B | 2,3,7,8-TCDF from both DB-5 and DB-225 | X | A |
| RSAI7-0.5B | All TCL compounds | X | A |
| RSAI7-0.5BDL | 2,3,7,8-TCDF from DB-5 | X | A |

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.

**Tronox LLC Facility, 2008 Phase B Investigation, Henderson, Nevada
Dioxins/Dibenzofurans - Data Qualification Summary - SDG E0800661**

| SDG | Sample | Compound | Flag | A or P | Reason (Code) |
|----------|----------------------------|--|--|--------|-------------------------------------|
| E0800661 | RSAJ8-0.5B | 2,3,7,8-TCDD 1,2,3,4,6,7,8-HpCDF | J+ (all detects) J+ (all detects) | P | Laboratory control samples (%R) (l) |
| E0800661 | RSAJ8-0.5B | OCDD OCDF | J (all detects) UJ (all non-detects) J (all detects) UJ (all non-detects) | P | Internal standards (%R) (i) |
| E0800661 | RSAI7-0.5B RSAI7-0.5BDL | 2,3,7,8-TCDF 1,2,3,4,7,8-HxCDD 1,2,3,6,7,8-HxCDD 1,2,3,7,8,9-HxCDD 1,2,3,4,6,7,8-HpCDD OCDD 1,2,3,4,7,8-HxCDF 1,2,3,6,7,8-HxCDF 2,3,4,6,7,8-HxCDF 1,2,3,7,8,9-HxCDF 1,2,3,4,6,7,8-HpCDF 1,2,3,4,7,8,9-HpCDF OCDF Total HxCDD Total HpCDD Total TCDF Total HxCDF Total HpCDF | J (all detects) UJ (all non-detects) | P | Internal standards (%R) (i) |
| E0800661 | RSAJ8-0.5B | 2,3,7,8-TCDF 1,2,3,4,7,8-HxCDF 1,2,3,4,6,7,8-HpCDF OCDF | J (all detects) J (all detects) J (all detects) J (all detects) | A | Project Quantitation Limit (e) |
| E0800661 | RSAI7-0.5B | 2,3,7,8-TCDD 1,2,3,7,8-PeCDD 1,2,3,4,7,8-HxCDD 1,2,3,6,7,8-HxCDD 1,2,3,7,8,9-HxCDD 1,2,3,4,6,7,8-HpCDD OCDD 2,3,7,8-TCDF 2,3,4,7,8-PeCDF 1,2,3,4,7,8-HxCDF 1,2,3,6,7,8-HxCDF 2,3,4,6,7,8-HxCDF 1,2,3,7,8,9-HxCDF 1,2,3,4,6,7,8-HpCDF 1,2,3,4,7,8,9-HpCDF OCDF | J (all detects) | A | Project Quantitation Limit (e) |

| SDG | Sample | Compound | Flag | A or P | Reason (Code) |
|----------|--|---|---|--------|---------------------------------|
| E0800661 | RSAI7-0.5BDL | 2,3,7,8-TCDF 1,2,3,4,7,8-HxCDF 1,2,3,4,6,7,8-HpCDF 1,2,3,4,7,8,9-HpCDF OCDF | J (all detects) J (all detects) J (all detects) J (all detects) J (all detects) | A | Project Quantitation Limit (e) |
| E0800661 | RSAJ8-0.5B RSAI7-0.5B RSAJ8-0.5BDL RSAI7-0.5BDL | All compounds reported below the PQL. | J (all detects) | A | Project Quantitation Limit (sp) |
| E0800661 | RSAJ8-0.5B RSAI7-0.5B RSAJ8-0.5BDL RSAI7-0.5BDL | All compounds reported as EMPC | JK (all detects) | A | Project Quantitation Limit (k) |
| E0800661 | RSAJ8-0.5B | 2,3,7,8-TCDF from both DB-5 and DB-225 | X | A | Overall assessment of data (o) |
| E0800661 | RSAI7-0.5B | All TCL compounds | X | A | Overall assessment of data (o) |
| E0800661 | RSAI7-0.5BDL | 2,3,7,8-TCDF from DB-5 | X | A | Overall assessment of data (o) |

**Tronox LLC Facility, 2008 Phase B Investigation, Henderson, Nevada
Dioxins/Dibenzofurans - Laboratory Blank Data Qualification Summary - SDG E0800661**

No Sample Data Qualified in this SDG

**Tronox LLC Facility, 2008 Phase B Investigation, Henderson, Nevada
Dioxins/Dibenzofurans - Field Blank Data Qualification Summary - SDG E0800661**

No Sample Data Qualified in this SDG

Tronox Northgate Henderson

VALIDATION COMPLETENESS WORKSHEET

LDC #: 21257U21

SDG #: E0800661

Laboratory: Columbia Analytical Services

Stage 2B

Date: 8/4/09

Page: 1 of 1

Reviewer: [Signature]

2nd Reviewer: [Signature]

METHOD: HRGC/HRMS Dioxins/Dibenzofurans (EPA SW 846 Method 8290)

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: 7/10-11/08 |
| II. | HRGC/HRMS instrument performance check | A | |
| III. | Initial calibration | A | |
| IV. | Routine calibration/ICV | AMW | |
| V. | Blanks | AMW | |
| VI. | Matrix spike/Matrix spike duplicates | N | |
| VII. | Laboratory control samples | AMW | LOS |
| VIII. | Regional quality assurance and quality control | N | |
| IX. | Internal standards | AMW | |
| X. | Target compound identifications | N | |
| XI. | Compound quantitation and CRQLs | AMW | |
| XII. | System performance | N | |
| XIII. | Overall assessment of data | AMW | |
| 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:

| | | | | | | | | |
|----------------|----|---------------|-----------------|--------------|-----------------|----------|----|---------|
| 1 | 1 | RSAJ8-0.5B | 11 ⁵ | EA0800294-01 | 21 ¹ | U129088 | 31 | U1290T2 |
| 2 ³ | 2 | RSAT7-0.5B | 12 | EA0800299-01 | 22 ¹ | C15311#2 | 32 | |
| 3 ² | 3 | RSAT8-0.5B DL | 13 | | 23 ² | C15322#2 | 33 | |
| 4 ⁴ | 4 | RSAT7-0.5B DL | 14 | | 24 ³ | U129088 | 34 | |
| 5 | 5 | | 15 | | 25 ^A | U217360 | 35 | |
| 6 | 6 | | 16 | | 26 ⁵ | U216848 | 36 | |
| 7 | 7 | | 17 | | 27 | | 37 | |
| 8 | 8 | | 18 | | 28 | | 38 | |
| 9 | 9 | | 19 | | 29 | | 39 | |
| 10 | 10 | | 20 | | 30 | | 40 | |

Notes: _____

VALIDATION FINDINGS WORKSHEET

METHOD: HRGC/HRMS Dioxins/Dibenzofurans (EPA SW 846 Method 8290)

| | | | | |
|----------------------|------------------------|------------------------|------------------------|----------------|
| A. 2,3,7,8-TCDD | F. 1,2,3,4,6,7,8-HpCDD | K. 1,2,3,4,7,8-HxCDF | P. 1,2,3,4,7,8,9-HpCDF | U. Total HpCDD |
| B. 1,2,3,7,8-PeCDD | G. OCDD | L. 1,2,3,6,7,8-HxCDF | Q. OCDF | V. Total TCDF |
| C. 1,2,3,4,7,8-HxCDD | H. 2,3,7,8-TCDF | M. 2,3,4,6,7,8-HxCDF | R. Total TCDD | W. Total PeCDF |
| D. 1,2,3,6,7,8-HxCDD | I. 1,2,3,7,8-PeCDF | N. 1,2,3,7,8,9-HxCDF | S. Total PeCDD | X. Total HxCDF |
| E. 1,2,3,7,8,9-HxCDD | J. 2,3,4,7,8-PeCDF | O. 1,2,3,4,6,7,8-HpCDF | T. Total HxCDD | Y. Total HpCDF |

Notes:

VALIDATION FINDINGS WORKSHEET
Blanks

LDC #: 2/257/2
 SDG #: See above

Page: 1 of 1
 Reviewer: _____
 2nd Reviewer: _____

METHOD: HRGC/HRMS Dioxins/Dibenzofurans (EPA Method 8290)

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

- N N/A Were all samples associated with a method blank?
- N N/A Was a method blank analyzed for each matrix?
- N N/A Was the blank contaminated? If yes, please see qualification below.

Blank extraction date: 7/14/08 **Blank analysis date:** 7/18/08

Conc. units: ng/L **Associated Samples:** 2, 4 (>5x)

| Compound | Blank ID | Sample Identification |
|----------|--------------|-----------------------|
| | 5A0800294-01 | |
| F | 0.265 | |
| G | 1.10 | |
| H | 0.581 | |
| I | 0.276 | |
| K | 0.340 | |
| L | 0.229 | |
| O | 0.405 | |

Blank extraction date: _____ **Blank analysis date:** _____
Conc. units: _____ **Associated Samples:** _____

| Compound | Blank ID | Sample Identification |
|----------|----------|-----------------------|
| | | |
| F | 0.177 | |
| R | 0.796 | |
| S | 0.0875 | |
| U | 0.5747 | |
| V | 2.213 | |
| W | 1.67 | |
| X | 1.14 | |

CIRCLED RESULTS WERE NOT QUALIFIED. ALL RESULTS NOT CIRCLED WERE QUALIFIED BY THE FOLLOWING STATEMENT:
 All contaminants within five times the method blank concentration were qualified as not detected, "U".

LDC #: 23516
 SDG #: Seecow

VALIDATION FINDINGS WORKSHEET
Overall Assessment of Data

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

METHOD: HRGC/HRMS Dioxins/Dibenzofurans (EPA SW 846 Method 8290)

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

All available information pertaining to the data were reviewed using professional judgement to complement the determination of the overall quality of the data.

N N/A Was the overall quality and usability of the data acceptable?

| # | Date | Sample ID | Finding | Associated Samples | Qualifications |
|---|------|-----------|----------------------|--------------------|----------------|
| 1 | | | H on PB-5 and DB-225 | 1 | X (07) |
| 2 | | | All | 2 | |
| 4 | | | H on PB-5 | 4 | |
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Comments: _____

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

Project/Site Name: Tronox LLC Facility, 2008 Phase B Investigation,
Henderson, Nevada

Collection Date: July 10, 2008

LDC Report Date: August 26, 2009

Matrix: Soil

Parameters: Dioxins/Dibenzofurans

Validation Level: Stage 2B

Laboratory: Columbia Analytical Services, Inc.

Sample Delivery Group (SDG): E0800662

Sample Identification

RSAL2-0.5B
RSAK2-0.5B*
RSAL2-0.5BDL
RSAL2-0.5BMS
RSAL2-0.5BMSD

*Indicates sample was analyzed for screening purposes only.

Introduction

This data review covers 5 soil samples listed on the cover sheet including dilutions and reanalysis as applicable. The analyses were per EPA SW 846 Method 8290 for Polychlorinated Dioxins/Dibenzofurans.

This review follows the Standard Operating Procedures (SOP) 40, Data Review/Validation (BRC 2009), the Quality Assurance Project Plan Tronox LLC Facility, Henderson, Nevada (June 2009), NDEP guidance (May 2006), and USEPA Contract Laboratory Program National Functional Guidelines for Polychlorinated Dioxins/Dibenzofurans Data Review (September 2005) as there are no current guidelines for the method 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.

Blank results are summarized in Section V.

Field duplicates are summarized in Section XIV.

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.
- B The analytical result may be a false positive totally attributable to blank contamination. This qualifier is applicable to radiochemistry analysis only.
- JB The analytical result may be biased high and partially attributable to blank contamination. This qualifier is applicable to radiochemistry analysis only.
- JK The analytical result is an estimated maximum possible concentration (EMPC).
- X The analytical result is not used for reporting because a more accurate and precise result is reported in its place.
- J-TDS The analytical result is estimated based on failure of the Total Dissolved Solids (TDS) correctness check performed in accordance with the Standard Method 1030E.
- J-CAB The analytical result is estimated based on failure of the cation-anion balance correctness check performed in accordance with Standard Method 1030E.
- J-TDS & CAB The analytical result is unreliable based on the failure of the cation-anion balance and TDS correctness check performed in accordance with standard Method 1030E.
- 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. HRGC/HRMS Instrument Performance Check

Instrument performance was checked at the required daily frequency.

Retention time windows were established for all homologues. The chromatographic resolution between 2,3,7,8-TCDD and peaks representing any other unlabeled TCDD isomer was less than or equal to 25%.

III. Initial Calibration

A five point initial calibration was performed as required by the method.

Percent relative standard deviations (%RSD) were less than or equal to 20.0% for unlabeled compounds and less than or equal to 30.0% for labeled compounds.

The ion abundance ratios for all PCDDs and PCDFs were within validation criteria.

IV. Routine Calibration (Continuing)

Routine calibration was performed at the required frequencies.

All of the routine calibration percent differences (%D) between the initial calibration RRF and the routine calibration RRF were less than or equal to 20.0% for unlabeled compounds and less than or equal to 30.0% for labeled compounds.

The ion abundance ratios for all PCDDs and PCDFs were within validation criteria.

V. Blanks

Method blanks were reviewed for each matrix as applicable. No polychlorinated dioxin/dibenzofuran contaminants were found in the method blanks with the following exceptions:

| Method Blank ID | Extraction Date | Compound | Concentration | Associated Samples |
|-----------------|-----------------|---------------------|---------------|--------------------|
| EQ0800294-01 | 7/14/08 | 1,2,3,4,6,7,8-HpCDD | 0.265 ng/Kg | RSAL2-0.5B |
| | | OCDD | 1.10 ng/Kg | RSAL2-0.5BDL |
| | | 2,3,7,8-TCDF | 0.581 ng/Kg | |
| | | 1,2,3,7,8-PeCDF | 0.276 ng/Kg | |
| | | 1,2,3,4,7,8-HxCDF | 0.340 ng/Kg | |
| | | 1,2,3,6,7,8-HxCDF | 0.229 ng/Kg | |
| | | 1,2,3,4,6,7,8-HpCDF | 0.405 ng/Kg | |
| | | 1,2,3,4,7,8,9-HpCDF | 0.177 ng/Kg | |
| | | OCDF | 0.796 ng/Kg | |
| | | Total PeCDD | 0.0845 ng/Kg | |
| | | Total HpCDD | 0.547 ng/Kg | |
| | | Total TCDF | 2.13 ng/Kg | |
| | | Total PeCDF | 1.67 ng/Kg | |
| | | Total HxCDF | 1.14 ng/Kg | |
| | | Total HpCDF | 0.603 ng/Kg | |

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

No field blanks were identified in this SDG.

VI. Matrix Spike/Matrix Spike Duplicates

Matrix spike (MS) and matrix spike duplicate (MSD) samples were reviewed for each matrix as applicable. Although the MS/MSD percent recoveries (%R) and relative percent differences (RPD) were not within QC limits for all compounds, the LCS percent recoveries (%R) were within QC limits and no data were qualified.

VII. Laboratory Control Samples (LCS)

Laboratory control samples were reviewed for each matrix as applicable. The percent recoveries (%R) were within the QC limits.

VIII. Regional Quality Assurance and Quality Control

Not applicable.

IX. Internal Standards

All internal standard recoveries were within QC limits with the following exceptions:

| Sample | Internal Standards | %R (Limits) | Compound | Flag | A or P |
|--------------|--|--|--|---|--------|
| RSAL2-0.5B | ¹³ C-1,2,3,4,7,8-HxCDF ¹³ C-1,2,3,6,7,8-HxCDD ¹³ C-1,2,3,4,6,7,8-HpCDF ¹³ C-1,2,3,4,6,7,8-HpCDD ¹³ C-OCDD | 22 (40-135) 24 (40-135) 16 (40-135) 14 (40-135) 8 (40-135) | 1,2,3,4,7,8-HxCDD 1,2,3,6,7,8-HxCDD 1,2,3,7,8,9-HxCDD 1,2,3,4,6,7,8-HpCDD OCDD 1,2,3,4,7,8-HxCDF 1,2,3,6,7,8-HxCDF 2,3,4,6,7,8-HxCDF 1,2,3,7,8,9-HxCDF 1,2,3,4,6,7,8-HpCDF 1,2,3,4,7,8,9-HpCDF OCDF Total HxCDD Total HpCDD Total HxCDF Total HpCDF | J (all detects) UJ (all non-detects) | P |
| RSAL2-0.5BDL | ¹³ C-1,2,3,4,7,8-HxCDF ¹³ C-1,2,3,6,7,8-HxCDD ¹³ C-1,2,3,4,6,7,8-HpCDF ¹³ C-1,2,3,4,6,7,8-HpCDD ¹³ C-OCDD | 24 (40-135) 29 (40-135) 21 (40-135) 17 (40-135) 7 (40-135) | 1,2,3,4,7,8-HxCDD 1,2,3,6,7,8-HxCDD 1,2,3,7,8,9-HxCDD 1,2,3,4,6,7,8-HpCDD OCDD 1,2,3,4,7,8-HxCDF 1,2,3,6,7,8-HxCDF 2,3,4,6,7,8-HxCDF 1,2,3,7,8,9-HxCDF 1,2,3,4,6,7,8-HpCDF 1,2,3,4,7,8,9-HpCDF OCDF Total HxCDD Total HpCDD Total HxCDF Total HpCDF | J (all detects) UJ (all non-detects) | P |

X. Target Compound Identifications

Raw data were not reviewed for this SDG.

XI. Project Quantitation Limit

All project quantitation limits were within validation criteria with the following exceptions:

| Sample | Compound | Finding | Criteria | Flag | A or P |
|--------------|--|---|---|--|--------|
| RSAL2-0.5B | 2,3,7,8-TCDD 1,2,3,7,8-PeCDD 1,2,3,4,7,8-HxCDD 1,2,3,6,7,8-HxCDD 1,2,3,7,8,9-HxCDD 1,2,3,4,6,7,8-HpCDD OCDD 2,3,7,8-TCDF 2,3,4,7,8-PeCDF 1,2,3,4,7,8-HxCDF 1,2,3,6,7,8-HxCDF 2,3,4,6,7,8-HxCDF 1,2,3,7,8,9-HxCDF 1,2,3,4,6,7,8-HpCDF 1,2,3,4,7,8,9-HpCDF OCDF | Sample result exceeded calibration range. | Reported result should be within calibration range. | J (all detects) J (all detects) J (all detects) J (all detects) J (all detects) J (all detects) J (all detects) J (all detects) J (all detects) J (all detects) J (all detects) J (all detects) J (all detects) J (all detects) J (all detects) J (all detects) | A |
| RSAL2-0.5BDL | OCDF | Sample result exceeded calibration range. | Reported result should be within calibration range. | J (all detects) | A |

All compounds reported below the PQL were qualified as follows:

| Sample | Finding | Flag | A or P |
|-----------------------------|---------------------------------------|-----------------|--------|
| All samples in SDG E0800662 | All compounds reported below the PQL. | J (all detects) | A |

All compounds reported as EMPC were qualified as follows:

| Sample | Finding | Flag | A or P |
|-----------------------------|--|------------------|--------|
| All samples in SDG E0800662 | All compounds reported as estimated maximum possible concentration (EMPC). | JK (all detects) | A |

Raw data were not reviewed for this SDG.

XII. System Performance

Raw data were not reviewed for this SDG.

XIII. Overall Assessment of Data

The overall assessment of data was acceptable. In the case where more than one result was reported for an individual sample, the least technically acceptable results were rejected as follows:

| Sample | Compound | Flag | A or P |
|--------------|------------------------|------|--------|
| RSAL2-0.5B | All TCL compounds | X | A |
| RSAL2-0.5BDL | 2,3,7,8-TCDF from DB-5 | X | A |

Sample RSAK2-0.5B was analyzed for screening purposes only.

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.

**Tronox LLC Facility, 2008 Phase B Investigation, Henderson, Nevada
Dioxins/Dibenzofurans - Data Qualification Summary - SDG E0800662**

| SDG | Sample | Compound | Flag | A or P | Reason (Code) |
|----------|----------------------------|--|---|--------|------------------------------------|
| E0800662 | RSAL2-0.5B RSAL2-0.5BDL | 1,2,3,4,7,8-HxCDD 1,2,3,6,7,8-HxCDD 1,2,3,7,8,9-HxCDD 1,2,3,4,6,7,8-HpCDD OCDD 1,2,3,4,7,8-HxCDF 1,2,3,6,7,8-HxCDF 2,3,4,6,7,8-HxCDF 1,2,3,7,8,9-HxCDF 1,2,3,4,6,7,8-HpCDF 1,2,3,4,7,8,9-HpCDF OCDF Total HxCDD Total HpCDD Total HxCDF Total HpCDF | J (all detects) UJ (all non-detects) | P | Internal standards (%R) (i) |
| E0800662 | RSAL2-0.5B | 2,3,7,8-TCDD 1,2,3,7,8-PeCDD 1,2,3,4,7,8-HxCDD 1,2,3,6,7,8-HxCDD 1,2,3,7,8,9-HxCDD 1,2,3,4,6,7,8-HpCDD OCDD 2,3,7,8-TCDF 2,3,4,7,8-PeCDF 1,2,3,4,7,8-HxCDF 1,2,3,6,7,8-HxCDF 2,3,4,6,7,8-HxCDF 1,2,3,7,8,9-HxCDF 1,2,3,4,6,7,8-HpCDF 1,2,3,4,7,8,9-HpCDF OCDF | J (all detects) J (all detects) J (all detects) J (all detects) J (all detects) J (all detects) J (all detects) J (all detects) J (all detects) J (all detects) J (all detects) J (all detects) J (all detects) J (all detects) J (all detects) J (all detects) J (all detects) | A | Project Quantitation Limit (e) |
| E0800662 | RSAL2-0.5BDL | OCDF | J (all detects) | A | Project Quantitation Limit (e) |
| E0800662 | RSAL2-0.5B RSAL2-0.5BDL | All compounds reported below the PQL. | J (all detects) | A | Project Quantitation Limit (sp) |
| E0800662 | RSAL2-0.5B RSAL2-0.5BDL | All compounds reported as EMPC | JK (all detects) | A | Project Quantitation Limit (k) |
| E0800662 | RSAL2-0.5B | All TCL compounds | X | A | Overall assessment of data (o) |
| E0800662 | RSAL2-0.5BDL | 2,3,7,8-TCDF from DB-5 | X | A | Overall assessment of data (o) |

**Tronox LLC Facility, 2008 Phase B Investigation, Henderson, Nevada
Dioxins/Dibenzofurans - Laboratory Blank Data Qualification Summary - SDG
E0800662**

No Sample Data Qualified in this SDG

**Tronox LLC Facility, 2008 Phase B Investigation, Henderson, Nevada
Dioxins/Dibenzofurans - Field Blank Data Qualification Summary - SDG E0800662**

No Sample Data Qualified in this SDG

Tronox Northgate Henderson

VALIDATION COMPLETENESS WORKSHEET

LDC #: 21257V21
 SDG #: E0800662
 Laboratory: Columbia Analytical Services

Stage 2B

Date: 8/14/09
 Page: 1 of 1
 Reviewer: [Signature]
 2nd Reviewer:

METHOD: HRGC/HRMS Dioxins/Dibenzofurans (EPA SW 846 Method 8290)

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: 7/10/08 |
| II. | HRGC/HRMS Instrument performance check | A | |
| III. | Initial calibration | A | |
| IV. | Routine calibration/ICV | A | |
| V. | Blanks | SW | |
| VI. | Matrix spike/Matrix spike duplicates | SW | |
| VII. | Laboratory control samples | SWA | LCS |
| VIII. | Regional quality assurance and quality control | N | |
| IX. | Internal standards | SW | |
| X. | Target compound identifications | N | |
| XI. | Compound quantitation and CRQLs | SW | |
| XII. | System performance | N | |
| XIII. | Overall assessment of data | SW | |
| XIV. | Field duplicates | N | |
| XV. | 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:

| | | | | | | |
|----|----------------------------------|----|------------------|----|---------|----|
| 1 | RSAL2-0.5B | 11 | ER0800294-01 | 21 | U216841 | 31 |
| 2 | RSAL2-0.5B RSAK2-0.5B | 12 | → screening only | 22 | U217360 | 32 |
| 3 | RSAL2-0.5BDC | 13 | | 23 | U216848 | 33 |
| 4 | RSAL2-0.5BMS | 14 | | 24 | | 34 |
| 5 | ↓ MS | 15 | | 25 | | 35 |
| 6 | | 16 | | 26 | | 36 |
| 7 | | 17 | | 27 | | 37 |
| 8 | | 18 | | 28 | | 38 |
| 9 | | 19 | | 29 | | 39 |
| 10 | | 20 | | 30 | | 40 |

Notes: # 2 is DIF-screen

VALIDATION FINDINGS WORKSHEET

METHOD: HRGC/HRMS Dioxins/Dibenzofurans (EPA SW 846 Method 8290)

| | | | | |
|----------------------|------------------------|------------------------|------------------------|----------------|
| A. 2,3,7,8-TCDD | F. 1,2,3,4,6,7,8-HpCDD | K. 1,2,3,4,7,8-HxCDF | P. 1,2,3,4,7,8,9-HpCDF | U. Total HpCDD |
| B. 1,2,3,7,8-PeCDD | G. OCDD | L. 1,2,3,6,7,8-HxCDF | Q. OCDF | V. Total TCDF |
| C. 1,2,3,4,7,8-HxCDD | H. 2,3,7,8-TCDF | M. 2,3,4,6,7,8-HxCDF | R. Total TCDD | W. Total PeCDF |
| D. 1,2,3,6,7,8-HxCDD | I. 1,2,3,7,8-PeCDF | N. 1,2,3,7,8,9-HxCDF | S. Total PeCDD | X. Total HxCDF |
| E. 1,2,3,7,8,9-HxCDD | J. 2,3,4,7,8-PeCDF | O. 1,2,3,4,6,7,8-HpCDF | T. Total HxCDD | Y. Total HpCDF |

Notes:

VALIDATION FINDINGS WORKSHEET
Blanks

METHOD: HRGC/HRMS Dioxins/Dibenzofurans (EPA Method 8290)

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

- Y N N/A Were all samples associated with a method blank?
- Y N N/A Was a method blank analyzed for each matrix?
- Y N N/A Was the blank contaminated? If yes, please see qualification below.

Blank extraction date: 7/14/08 Blank analysis date: 7/18/08

Conc. units: ng/kg Associated Samples: 1, 3 (>5x)

| Compound | Blank ID | Sample Identification |
|--------------|------------------|-----------------------|
| F | 0.265 | |
| G | 1.10 | |
| H | 0.581 | |
| I | 0.276 | |
| K | 0.340 | |
| L | 0.229 | |
| O | 0.405 | |

Blank extraction date: _____ Blank analysis date: _____
 Conc. units: _____ Associated Samples: _____

| Compound | Blank ID | Sample Identification |
|--------------|-------------------|-----------------------|
| P | 0.177 | |
| R | 0.796 | |
| S | 0.0845 | |
| U | 0.547 | |
| V | 0.213 | |
| W | 1.67 | |
| X | 1.14 | |
| Y | 0.603 | |

CIRCLED RESULTS WERE NOT QUALIFIED. ALL RESULTS NOT CIRCLED WERE QUALIFIED BY THE FOLLOWING STATEMENT:
 All contaminants within five times the method blank concentration were qualified as not detected, "U".

VALIDATION FINDINGS WORKSHEET
Matrix Spike/Matrix Spike Duplicates

LDC #: 21-257121
 SDG #: See cover

Page: 1 of 1
 Reviewer: g
 2nd Reviewer: _____

METHOD: HRGC/HRMS Dioxins/Dibenzofurans (EPA SW 846 Method 8290)

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

Y Were a matrix spike (MS) and matrix spike duplicate (MSD) analyzed for each matrix in this SDG? If no, indicate which matrix does not have an associated MS/MSD, Soil / Water.
N N/A
Y N/A
N N/A

Y Were the MS/MSD percent recoveries (%R) and the relative percent differences (RPD) within the QC limits?
Y N/A
N N/A

| # | Date | MS/MSD ID | Compound | MS %R (Limits) | MSD %R (Limits) | RPD (Limits) | Associated Samples | Qualifications |
|---|------|------------|--|----------------|-----------------|--------------|--------------------|----------------|
| | | <u>4/5</u> | <u>All DR and RPD out due to event</u> | | | | <u>1</u> | <u>No Qual</u> |
| | | | <u>> 4 X SA, LGS in</u> | () | () | () | | |
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VALIDATION FINDINGS WORKSHEET
Internal Standards

METHOD: HRGC/HRMS Dioxins/Dibenzofurans (EPA SW 846 Method 8290)
 Please see qualifications below for all questions answered "N". Not applicable questions are identified as "N/A".
 Are all internal standard recoveries within the 40-135% criteria?
 Y/N N/A
 Was the S/N ratio all internal standard peaks ≥ 10 ?
 Y/N N/A

Reason code (I)

| # | Date | Lab ID/Reference | Internal Standard | % Recovery (Limit: 40-135%) | Qualifications |
|---|------|------------------|-------------------|-----------------------------|---------------------------|
| | | 1 | Z | 22 | (40-135) ✓ N/A (C-F, K-X) |
| | | | F | 24 | () ✓ T-U, X-Y |
| | | | G | 16 | () |
| | | | H | 12 | () |
| | | | J | 8 | () |
| | | | | | |
| | | 3 | Z | 24 | () |
| | | | F | 29 | () |
| | | | G | 21 | () |
| | | | H | 17 | () |
| | | | I | 7 | () |
| | | | | | |
| | | 4 (M2) | A | 39 | () No Qual |
| | | | Z | 21 | () |
| | | | H | 22 | () |
| | | | S | 10 | () |
| | | | H | 14 | () |
| | | | I | 7 | () |
| | | | | | |
| | | | | | |

| Internal Standards | Check Standard Used | Internal Standards | Check Standard Used |
|--|---------------------|--------------------------------------|---------------------|
| A. ¹³ C-2,3,7,8-TCDF | | I. ¹³ C-OCDD | |
| B. ¹³ C-2,3,7,8-TCDD | | K. ¹³ C-1,2,3,4-TCDD | |
| C. ¹³ C-1,2,3,7,8-PeCDF | | L. ¹³ C-1,2,3,7,8,9-HxCDD | |
| D. ¹³ C-1,2,3,7,8-PeCDD | | M. | |
| E. ¹³ C-1,2,3,4,7,8-HxCDF | | N. | |
| F. ¹³ C-1,2,3,6,7,8-HxCDD | | O. | |
| G. ¹³ C-1,2,3,4,6,7,8-HpCDF | | P. | |
| H. ¹³ C-1,2,3,4,6,7,8-HpCDD | | | |

VALIDATION FINDINGS WORKSHEET

Internal Standards

METHOD: HRGC/HRMS Dioxins/Dibenzofurans (EPA SW 846 Method 8290)

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

~~Y/N~~ N/A Are all internal standard recoveries were within the 40-135% criteria?

~~Y/N~~ N/A Was the S/N ratio all internal standard peaks > 10 ?

| # | Date | Lab ID/Reference | Internal Standard | % Recovery (Limit: 40-135%) | Qualifications |
|---|------|------------------|-------------------|-----------------------------|----------------|
| | | 5 (MSD) | A | 33 (40-135) | No Qual |
| | | | B | 20 | |
| | | | F | 21 | |
| | | | G | 11 | |
| | | | H | 12 | |
| | | | I | 9 | |
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| | Internal Standards | Check Standard Used | Internal Standards | Check Standard Used |
|----|-------------------------------------|---------------------|--------------------|-----------------------------------|
| A. | ¹³ C-2,3,7,8-TCDF | | I. | ¹³ C-OCDD |
| B. | ¹³ C-2,3,7,8-TCDD | | K. | ¹³ C-1,2,3,4-TCDD |
| C. | ¹³ C-1,2,3,7,8-PeCDF | | L. | ¹³ C-1,2,3,7,8,9-HxCDD |
| D. | ¹³ C-1,2,3,7,8-PeCDD | | M. | |
| E. | ¹³ C-1,2,3,4,7,8-HxCDF | | N. | |
| F. | ¹³ C-1,2,3,6,7,8-HxCDD | | O. | |
| G. | ¹³ C-1,2,3,4,6,7,8-HpCDF | | P. | |
| H. | ¹³ C-1,2,3,4,6,7,8-HpCDD | | | |

VALIDATION FINDINGS WORKSHEET
Compound Quantitation and Reported CRQLs

LDC #: 21257121
 SDG #: [Signature]

METHOD: HRGC/HRMS Dioxins/Dibenzofurans (EPA SW 846 Method 8290)

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

Y N / N/A
 Y N / N/A

Were the correct internal standard (IS), quantitation ions and relative response factors (RRF) used to quantitate the compound? Compound quantitation and CRQLs were adjusted to reflect all sample dilutions and dry weight factors (if necessary).

| # | Date | Sample ID | Finding | Associated Samples | Qualifications |
|---|------|-----------|---------------------------------------|--------------------|------------------|
| 1 | | | <i>cpds > calib range</i> A → B | 1 | <i>plots / A</i> |
| 3 | | | B | 3 | <i>↓</i> |
| | | | | | |
| | | <i>AM</i> | <i>supc results</i> | <i>AM</i> | <i>JK (k)</i> |
| | | | | | |
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Comments: See sample calculation verification worksheet for recalculations

VALIDATION FINDINGS WORKSHEET
 Overall Assessment of Data

METHOD: HRGC/HRMS Dioxins/Dibenzofurans (EPA SW 846 Method 8290)

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

All available information pertaining to the data were reviewed using professional judgement to compliment the determination of the overall quality of the data.

Y N N/A Was the overall quality and usability of the data acceptable?

| # | Date | Sample ID | Finding | Associated Samples | Qualifications |
|---|------|-----------|---------------------|--------------------|----------------|
| 1 | | | All COB-5 and DB-25 | 1 | X (07) |
| 3 | | | H on DB-5 | 3 | ↓ |
| 2 | | | Sampling data only | 2 | Text |
| | | | | | |
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Comments: _____

