

**2006 SOURCE TEST REPORT**

**EMISSION MONITORING OF TWO LANDFILL GAS-FIRED  
FLARES AT THE REDWOOD LANDFILL**

**LANDFILL GAS CONTROL - FLARES - SOURCES A-50 AND A-51  
FACILITY NUMBER A1179**

Test dates: June 7 and 8, 2006

Submittal Date: July 31, 2006

Prepared for:

Waste Management, Inc.  
8950 Redwood Highway  
Novato, CA 94948

Prepared by:



Shaw Environmental, Inc.  
2360 Bering Drive  
San Jose, California 95131

**Project No.: 121105.01**

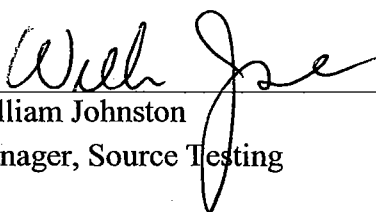
## **2006 SOURCE TEST REPORT**

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The material and data in this report were reviewed by the undersigned.



  
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William Johnston  
Manager, Source Testing

## SUMMARY OF 2006 SOURCE TEST INFORMATION

Source & Location: Enclosed Gas Flares A-50 and A-51  
Waste Management, Inc.  
8950 Redwood Highway  
Novato, CA 94948

Site Contact: Beth Shiverdecker: 415-892-2851

Source Description: A-50: Power Strategies EV 4000 enclosed ground flare assembly  
with a total capacity of 4000 scfm  
A-51: Perennial Energy Inc., 90 MMBtu/hr, enclosed ground flare  
assembly with a total capacity of 3000 scfm

Reference: Major Facility Review Permit  
Condition Number 19867  
Emission Limits:  
CO = 0.30 lb/MMBtu (123 ppmv at 15% O<sub>2</sub>)  
NO<sub>x</sub> = 0.6 lb/MMBtu (15 ppmv at 15% O<sub>2</sub>)  
NMOC = 30 ppmv at 3% O<sub>2</sub>

Agency: Bay Area Air Quality Management District  
939 Ellis Street  
San Francisco, CA 94109

Source Test Contractor: Shaw Environmental, Inc  
2360 Bering Drive  
San Jose, CA 95131

Contact: William Johnston: 408-382-5800

Test Dates: June 7 and 8, 2006, for A-50 and A-51, respectively

### Notes:

scfm = standard cubic feet per minute  
MMBtu/hr = million British thermal units per hour  
CO = carbon monoxide  
lb/MMBtu = pounds per million British thermal units  
ppmv = parts per million by volume  
O<sub>2</sub> = oxygen  
NO<sub>x</sub> = oxides of nitrogen  
NMOC = non-methane organic compounds

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## 1.0 Introduction

### 1.1 Summary

Shaw Environmental, Inc. (Shaw) performed source tests of landfill gas (LFG) Flares A-50 and A-51 at the Redwood Landfill located in Novato, California. This report presents the results of the test program. Tables 1 and 2 summarize the test results for Flares A-50 and A-51, respectively, as compared to the emission limits. The flares met all compliance criteria.

**Table 1. Compliance Summary  
Redwood Landfill Flare A-50  
Test date: June 7, 2006**

| Condition                                        | Average Result | Emission Limit | Compliance Status |
|--------------------------------------------------|----------------|----------------|-------------------|
| NOx (ppmv @ 15% O <sub>2</sub> )                 | 11             | 15             | In Compliance     |
| CO (ppmv @ 15% O <sub>2</sub> )                  | 35             | 123            | In Compliance     |
| NMOC (ppmv @ 3% O <sub>2</sub> , dry as methane) | 4              | 30             | In Compliance     |

*Notes:*

NOx = oxides of nitrogen

CO = carbon monoxide

NMOC = non-methane organic compounds

ppmv = parts per million by volume

O<sub>2</sub> = oxygen

**Table 2. Compliance Summary  
Redwood Landfill Flare A-51  
Test date: June 8, 2006**

| Condition                                        | Average Result | Emission Limit | Compliance Status |
|--------------------------------------------------|----------------|----------------|-------------------|
| NOx (ppmv @ 15% O <sub>2</sub> )                 | 13             | 15             | In Compliance     |
| CO (ppmv @ 15% O <sub>2</sub> )                  | 12             | 123            | In Compliance     |
| NMOC (ppmv @ 3% O <sub>2</sub> , dry as methane) | 7              | 30             | In Compliance     |

## **1.2 Overview**

Shaw, an approved California Air Resources Board (ARB) source test contractor, performed the source test program. The test program was designed to demonstrate compliance with the Bay Area Air Quality Management District (BAAQMD) Major Facility Review Permit (MFRP) A1179, Condition Number 19867. The source test was performed on the flare exhaust to assess the emissions of NMOC, NO<sub>x</sub>, and CO. Additionally, an LFG sample was collected for gas characterization.

The source tests were conducted on June 7 and 8, 2006 for Flares A-50 and A-51, respectively. Sampling was performed by Mr. Bill Johnston and Mr. Scott Borderieux of Shaw. The BAAQMD was notified of the test via a fax dated May 24, 2006, that included the Source Test Protocol. The BAAQMD did not have a representative present for the test. The source test program for collecting the required performance data is discussed in Section 2; the source testing results are presented in Section 3; the source test procedures are presented in Section 4; and data review is discussed in Section 5. Source test data are located in Appendix A. Continuous emission monitoring data are located in Appendix B. Process data are located in Appendix C. LFG analytical results are presented in Appendix D. Copies of the Source Test Protocol and of Shaw's ARB Independent Contractor Program certification are presented in Appendix E. Redwood Landfill's MFRP is located in Appendix F. A site picture is located in Appendix G.

## **1.3 Source Operating Conditions**

### **1.3.1 Source A-50**

The A-50 flare is a Power Strategies EV 4000 enclosed ground flare assembly with a total capacity of 4000 standard cubic feet per minute (scfm). The flare is utilized as a back-up flare to flare A-51. The LFG volumetric flow rate is measured with a LFG mass flow meter. Three thermocouples are installed at the top, middle, and bottom of the flare exhaust stack to provide temperature indication for control of the combustion temperature. The flare was set at an operating temperature of 1,650 degrees Fahrenheit (°F) measured at the top thermocouple for the source test. The average flare temperature was 1,647 °F with an average LFG flow rate of 872 scfm. The LFG landfill gas flow rate and flare operating temperature are continuously recorded on a data acquisition system (DAS) and circular paper chart recorder.

Process data documenting the operation of the flare are contained in Appendix E. A picture of the flares is presented in Appendix G.

### **1.3.2 Source A-51**

The A-51 flare is a Perennial Energy Inc., 90 million British thermal units per hour (MMBtu/hr), enclosed ground flare assembly with a total capacity of 3,000 scfm. The LFG volumetric flow rate is measured with a mass flow meter. Three thermocouples are located at the top, middle, and bottom of the flare exhaust stack to provide temperature indication for control of the combustion temperature. The flare was set at an operating temperature of 1,450 °F measured at the top thermocouple for the source test. The average flare temperature was 1,450 °F with an average LFG flow rate of 2,603 scfm. The LFG landfill gas flow rate and flare operating temperature are continuously recorded on a DAS and strip paper chart recorder.

## 2.0 Source Test Program

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The objective of the flare source tests was to demonstrate compliance with requirements of MFRP Condition Number 19867 for LFG parameters NO<sub>x</sub>, CO, NMOC, and other gaseous constituents.

Measurements of each flare's emissions were conducted at the flare exhausts. LFG was sampled at the inlet header. Table 3 provides a matrix of the source test methods used at each sample location. U.S. Environmental Protection Agency (EPA), American Society for Testing and Materials (ASTM) and BAAQMD test methods were used. The source test program consisted of three 30-minute runs to measure gaseous emissions in each flare exhaust and one 30-minute run for the inlet header. The continuous emission monitoring system (CEMS) equipment was leak-checked before sampling began. Each instrument was calibrated before and after each sampling run with EPA protocol calibrated gas standards. The CEMS data are presented in Appendix B.

Concurrent with each flare test, Shaw collected one integrated 1-liter Tedlar<sup>®</sup> bag sample of the inlet LFG for analysis of NMOC, British thermal units (Btu), F-Factor, and additional fixed gases. The analytical laboratory results are contained in Appendix D.

The flare operating temperature and LFG flow rate were manually recorded on a field data sheet for each run. The outlet volumetric flow rate was calculated using the fuel flow rate, Btu, exhaust gas O<sub>2</sub> content, and the F-Factor according to EPA Method 19.

Results of the tests for the LFG and flare exhaust are summarized in Section 3. Compliance with the NMOC limit was demonstrated by verifying that the concentrations of NMOC were less than 30 ppmv at 3 percent O<sub>2</sub> as methane. The source test demonstrated that the flares operate with criteria pollutant emissions below the BAAQMD emission limits.

**Table 3. Source Test Methods and Instrumentation**

| Method                                                     | Inlet (fuel) Analyte                       | Exhaust Analyte           | Instrumentation/<br>Laboratory             | Principle                                             |
|------------------------------------------------------------|--------------------------------------------|---------------------------|--------------------------------------------|-------------------------------------------------------|
| ASTM D-1945/D-3588,<br>EPA Method 25 C<br>EPA Method TO 15 | Fixed Gases, Btu, F-Factor<br>NMOC<br>VOCs | —                         | Atmospheric Analysis &<br>Consulting, Inc. | Chromatographic Analysis<br>GC/FID/TCA Analysis<br>GC |
| BAAQMD Method ST-6                                         | —                                          | CO                        | Horiba VIA 510                             | IR                                                    |
| BAAQMD Method ST-7                                         |                                            | THC/CH <sub>4</sub> /NMOC | TECO Model 51                              | FID                                                   |
| BAAQMD Method ST-13A                                       |                                            | NO <sub>x</sub>           | Ecophysics Model 70E                       | Chemiluminescence                                     |
| BAAQMD Method ST-14                                        |                                            | O <sub>2</sub>            | Teledyne                                   | Fuel Cell                                             |
| EPA Method 19                                              |                                            | Flow rate, dscfm          | —                                          | Algorithm                                             |

*Notes:*

*VOCs = volatile organic compounds*

*THC = total hydrocarbons*

*CH<sub>4</sub> = methane*

*dscfm = dry standard cubic feet per minute*

*FID = flame ionization detection*

*GC = gas chromatography*

*TCA = thermal conductivity analyzer*

*IR = infrared*

### 3.0 Source Test Results

The source tests were performed on the LFG fuel and at the flare exhausts. Tables 4 and 5 present the results of the LFG fuel analysis and flare operating parameters as recorded by the DAS for Flares A-50 and A-51, respectively. One LFG sample was drawn per flare test for analysis of Btu, F-Factor, and additional fixed gases. Tables 6 and 7 present the results of the source tests for Flares A-50 and A-51, respectively. Triplicate 30-minute runs were performed on the flare exhaust to assess the NMOC, NO<sub>x</sub>, and CO emissions. Concurrent with each run, the O<sub>2</sub> content of the exhaust gas was also measured. Concentrations of CO and NO<sub>x</sub> are reported in units of ppmv corrected to 15 percent O<sub>2</sub>. Concentrations of NMOC are reported in units of ppmv corrected to 3 percent O<sub>2</sub>.

One LFG sample was analyzed for compliance with Condition No. 19867, Parts 18 and 31 (the LFG gas characterization).

**Table 4. Redwood Landfill  
Flare A-50, Operating Conditions and LFG Analysis  
Test date: June 7, 2006**

|                                      | Run 1     | Run 2     | Run 3     | Average |
|--------------------------------------|-----------|-----------|-----------|---------|
| Time                                 | 0811-0841 | 0852-0922 | 0931-1001 | —       |
| Flare Temperature (°F)               | 1,647     | 1,647     | 1,648     | 1,647   |
| Recorded Fuel Flow Rate (scfm)       | 874       | 872       | 870       | 872     |
| Heat Input (MMBtu/day)               | 653       | 652       | 650       | 652     |
| O <sub>2</sub> (%) <sup>1</sup>      | —         | 2.5       | —         | —       |
| CO <sub>2</sub> (%) <sup>1</sup>     | —         | 34        | —         | —       |
| N <sub>2</sub> (%) <sup>1</sup>      | —         | 13        | —         | —       |
| Methane (%) <sup>1</sup>             | —         | 51.2      | —         | —       |
| H <sub>2</sub> S (ppmv) <sup>1</sup> | —         | 92        | —         | —       |
| NMOC (ppmv) <sup>1</sup>             | —         | 456       | —         | —       |

<sup>1</sup> Only one sample of landfill gas was collected and analyzed for this value

CO<sub>2</sub> = carbon dioxide

N<sub>2</sub> = nitrogen

H<sub>2</sub>S = hydrogen sulfide

**Table 5. Redwood Landfill**  
**Flare A-51, Operating Conditions and LFG Analysis**  
**Test date: June 8, 2006**

|                                      | Run 1     | Run 2     | Run 3     | Average |
|--------------------------------------|-----------|-----------|-----------|---------|
| Time                                 | 0852-0922 | 0933-1003 | 1014-1044 | —       |
| Flare Temperature (°F)               | 1,448     | 1,452     | 1,451     | 1,450   |
| Recorded Fuel Flow Rate (scfm)       | 2,618     | 2,601     | 2,591     | 2,603   |
| Heat Input (MMBtu/day)               | 1,866     | 1,854     | 1,847     | 1,856   |
| O <sub>2</sub> (%) <sup>1</sup>      | —         | 3.2       | —         | —       |
| CO <sub>2</sub> (%) <sup>1</sup>     | —         | 33        | —         | —       |
| N <sub>2</sub> (%) <sup>1</sup>      | —         | 15        | —         | —       |
| Methane (%) <sup>1</sup>             | —         | 48.8      | —         | —       |
| H <sub>2</sub> S (ppmv) <sup>1</sup> | —         | 327       | —         | —       |
| NMOC (ppmv) <sup>1</sup>             | —         | 213       | —         | —       |

<sup>1</sup> Only one sample of landfill gas was collected and analyzed for this value.

**Table 6. Redwood Landfill**  
**Source Test Results — Landfill Flare Exhaust A-50**  
**Test date: June 7, 2006**

|                                              | Run 1     | Run 2     | Run 3     | Average | Emission Limit  |
|----------------------------------------------|-----------|-----------|-----------|---------|-----------------|
| Time                                         | 0811-0841 | 0852-0922 | 0931-1001 | —       | —               |
| Fuel Factor                                  | 9,147     | 9,147     | 9,147     | 9,147   | —               |
| Flare Temperature (°F)                       | 1,647     | 1,647     | 1,648     | 1,647   | —               |
| Landfill Gas Fuel Flow Rate (scfm)           | 874       | 872       | 870       | 872     | —               |
| Exhaust Flow Rate (dscfm)                    | 8,849     | 8,739     | 8,899     | 8,829   | —               |
| O <sub>2</sub> (%)                           | 11.1      | 11.0      | 11.2      | 11.1    | —               |
| <b>NO<sub>x</sub></b>                        |           |           |           |         |                 |
| NO <sub>x</sub> (ppmv)                       | 18.7      | 18.3      | 18.1      | 18.4    | —               |
| NO <sub>x</sub> (ppmv @ 15% O <sub>2</sub> ) | 11.3      | 10.9      | 11.0      | 11.1    | 15              |
| NO <sub>x</sub> (lb/hr)                      | 1.2       | 1.1       | 1.1       | 1.2     | —               |
| NO <sub>x</sub> (lb/MMBtu)                   | 0.043     | 0.042     | 0.042     | 0.043   | 0.06            |
| <b>CO</b>                                    |           |           |           |         |                 |
| CO (ppmv)                                    | 53.7      | 53.6      | 65.0      | 57.4    | —               |
| CO (ppmv @ 15% O <sub>2</sub> )              | 32.3      | 31.9      | 39.5      | 34.6    | 123             |
| CO (lb/hr)                                   | 2.1       | 2.0       | 2.5       | 2.2     | —               |
| CO (lb/MMBtu)                                | 0.076     | 0.075     | 0.093     | 0.081   | 0.30            |
| <b>Hydrocarbons as Methane</b>               |           |           |           |         |                 |
| NMOC (ppmv)                                  | 1.1       | 1.8       | 3.7       | 2.2     | —               |
| NMOC (ppmv corrected to 3% O <sub>2</sub> )  | 2.0       | 3.3       | 6.8       | 4.0     | 30              |
| NMOC Removal Efficiency (%) <sup>1</sup>     | —         | 96.0      | —         | —       | 98 <sup>2</sup> |

<sup>1</sup> Only one sample of landfill gas was analyzed to calculate this value.

<sup>2</sup> The efficiency requirement is utilized only if the outlet NMOC concentration is greater than 30 ppmv @ 3% oxygen.

**Table 7. Redwood Landfill**  
**Source Test Results — Landfill Flare Exhaust A-51**  
**Test date: June 8, 2006**

|                                              | Run 1     | Run 2     | Run 3     | Average | Emission Limit  |
|----------------------------------------------|-----------|-----------|-----------|---------|-----------------|
| Time                                         | 0852-0922 | 0933-1003 | 1014-1044 | —       | —               |
| Fuel Factor                                  | 9,158     | 9,158     | 9,158     | 9,158   | —               |
| Flare Temperature (°F)                       | 1,448     | 1,452     | 1,451     | 1,450   | —               |
| Landfill Gas Fuel Flow Rate (scfm)           | 2,618     | 2,601     | 2,591     | 2,603   | —               |
| Exhaust Flow Rate (dscfm)                    | 31,397    | 30,804    | 32,300    | 31,500  | —               |
| O <sub>2</sub> (%)                           | 13.0      | 12.9      | 13.3      | 13.1    | —               |
| <b>NO<sub>x</sub></b>                        |           |           |           |         |                 |
| NO <sub>x</sub> (ppmv)                       | 17.0      | 17.2      | 15.7      | 16.6    | —               |
| NO <sub>x</sub> (ppmv @ 15% O <sub>2</sub> ) | 12.7      | 12.7      | 12.2      | 12.5    | 15              |
| NO <sub>x</sub> (lb/hr)                      | 3.8       | 3.8       | 3.6       | 3.7     | —               |
| NO <sub>x</sub> (lb/MMBtu)                   | 0.049     | 0.049     | 0.047     | 0.048   | 0.06            |
| <b>CO</b>                                    |           |           |           |         |                 |
| CO (ppmv)                                    | 13.7      | 11.6      | 21.3      | 15.5    | —               |
| CO (ppmv @ 15% O <sub>2</sub> )              | 10.2      | 8.6       | 16.5      | 11.8    | 123             |
| CO (lb/hr)                                   | 1.9       | 1.6       | 3.0       | 2.1     | —               |
| CO (lb/MMBtu)                                | 0.024     | 0.020     | 0.039     | 0.028   | 0.30            |
| <b>Hydrocarbons as Methane</b>               |           |           |           |         |                 |
| NMOC (ppmv)                                  | 2.3       | 1.3       | 5.0       | 2.9     | —               |
| NMOC (ppmv corrected to 3% O <sub>2</sub> )  | 5.2       | 2.9       | 11.8      | 6.6     | 30              |
| NMOC Removal Efficiency (%) <sup>1</sup>     | —         | 92.8      | —         | —       | 98 <sup>2</sup> |

<sup>1</sup> Only one sample of landfill gas was analyzed to calculate this value.

<sup>2</sup> The efficiency requirement is utilized only if the outlet NMOC concentration is greater than 30 ppmv @ 3% oxygen.

**Table 8. Summary of Results, Redwood Landfill**  
**Landfill Gas Analysis**  
**Test date: June 8, 2006**

| Compound                                 | Measured Concentration (ppbv) | Permit Concentration Limit (ppbv) |
|------------------------------------------|-------------------------------|-----------------------------------|
| <b>Organic Compounds<sup>1</sup></b>     |                               |                                   |
| Acrylonitrile                            | < 1.0                         | 280                               |
| Benzene                                  | 328                           | 340                               |
| Carbon Tetrachloride                     | < 1.0                         | 70                                |
| Chlorobenzene                            | < 1.0                         | —                                 |
| Chlorodifluoromethane                    | 575                           | —                                 |
| Chloroethane                             | < 1.0                         | —                                 |
| Chloroform                               | < 1.0                         | 70                                |
| 1,1-Dichloroethane                       | < 1.0                         | 150                               |
| 1,1-Dichloroethene                       | < 1.0                         | —                                 |
| 1,2-Dichloroethane                       | < 1.0                         | 70                                |
| 1,4-Dichlorobenzene                      | < 1.0                         | 400                               |
| Dichlorodifluoromethane                  | 642                           | —                                 |
| Dichlorofluoromethane                    | < 1.0                         | —                                 |
| Ethylbenzene                             | 1,790                         | —                                 |
| 1,2-Dibromomethane                       | < 1.0                         | 70                                |
| Fluorotrichloromethane                   | < 1.0                         | —                                 |
| Hexane                                   | 471                           | —                                 |
| Isopropanol                              | 269                           | —                                 |
| Methyl Ethyl Ketone                      | 1,200                         | —                                 |
| Methylene Chloride                       | < 1.0                         | 320                               |
| Perchloroethylene (Tetrachloroethylene)  | < 1.0                         | 450                               |
| Toluene                                  | 4,200                         | —                                 |
| 1,1,1-Trichloroethane                    | < 1.0                         | —                                 |
| 1,1,2,2-Tetrachloroethane                | < 1.0                         | 70                                |
| Trichloroethylene                        | < 1.0                         | 250                               |
| Vinyl Chloride                           | 431                           | 880                               |
| Xylenes                                  | 2500                          | —                                 |
| Compound                                 | Measured Concentration (ppmv) | Permit Concentration Limit (ppmv) |
| <b>Sulfur Compounds<sup>1</sup></b>      |                               |                                   |
| Carbon Disulfide                         | < 0.05                        | —                                 |
| Carbonyl Sulfide                         | < 0.05                        | —                                 |
| Dimethyl Sulfide                         | < 0.05                        | —                                 |
| Ethyl Mercaptan                          | < 0.05                        | —                                 |
| Total Reduced Sulfur as Hydrogen Sulfide | 327                           | 1,300                             |
| Methyl Mercaptan                         | < 0.05                        | —                                 |
| <b>NMOC</b>                              |                               |                                   |
| NMOC (as Hexane)                         | 76                            | 750                               |

Notes:

ppbv = parts per billion by volume

ppmv = parts per million by volume

<sup>1</sup>For compounds reported as non-detected, a less than (<) sign is used and the method reporting limit value is reported.

## 4.0 Test Procedures

Shaw conducted the source test according to ASTM, BAAQMD and EPA test methods as described below.

### 4.1 Continuous Emissions Monitoring

Shaw measured stack gas constituent concentrations, using its CEMS sampling van, according to BAAQMD methods for NMOC, NO<sub>x</sub>, CO, and O<sub>2</sub>. Figure 1 is a schematic diagram of the Shaw CEMS.

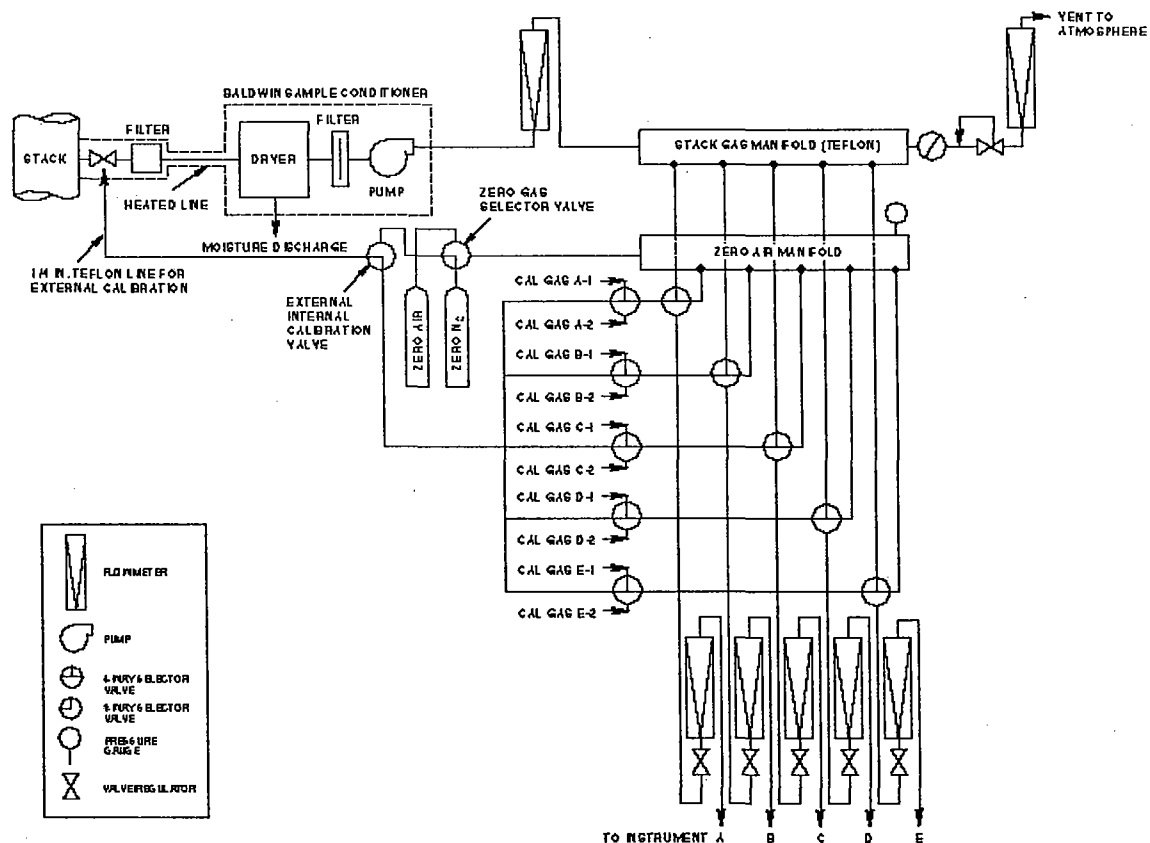


Figure 1. Schematic Diagram of the Shaw CEMS

In this source test, a sample pump was used to extract the sample gas continuously from each source through a stainless-steel sampling probe equipped with an in-stack filter, followed by a heated Teflon™ sample line and a Baldwin sample conditioner. The sample line between the sample point and the sample conditioner was heated to 250°F to maintain the gas stream temperatures above the water dew point. Gas flow was controlled with a back-pressure regulator. The sample flow rate to each instrument was controlled with individual rotameters. Instrument output was recorded with a strip chart recorder and a digital data logger, the STRATA data acquisition system. Table 6 lists the specifications of the instruments used in the source test.

**Table 9. CEMS Instrument Specifications**

| Gas             | Analyzer/Manufacturer | Principle of Operation | Operating Range              |
|-----------------|-----------------------|------------------------|------------------------------|
| O <sub>2</sub>  | Teledyne 326A         | Fuel Cell              | 0-25%                        |
| THC             | TECO Model 51         | Flame Ionization       | 0-50 ppmv as CH <sub>4</sub> |
| NO <sub>x</sub> | Ecophysics 70E        | Chemiluminescence      | 0-50 ppmv                    |
| CO              | Horiba VIA 510        | NDIR                   | 0-100 ppmv                   |

Note:

NDIR = nondispersive infrared

Before the start of testing, the sampling system was assembled at the site and leak-checked. Calibrations were performed on each of the instruments using certified calibration gases.

A span drift check was performed at the end of each run to measure analyzer drift. The drift check measurement uses a zero gas (N<sub>2</sub>) and a high calibration gas. All of the instrument checks conducted for this source test were within the drift limit criterion of 3 percent, as specified in the test methods. Reported values include a correction for the recorded instrument drift, as specified in EPA Method 6C. Copies of data logger printouts, calibration data sheets, strip charts, and gas certificates of analysis, are presented in Appendix B.

## 4.2 Landfill Gas

Shaw measured LFG constituent concentrations according to EPA and ASTM test methods. The sample was collected from a sample port located at the inlet header. The sample was collected in a 1-liter Tedlar® bag through Teflon™ tubing. The sample line was purged with sample gas prior to sampling. The sample flow rate was controlled with rotameters to collect a 30-minute integrated sample. The sample was kept out of sunlight and analyzed within the method holding times. Analytical results are presented in Appendix D.

## 5.0 Data Review

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The Air Quality Group within Shaw performed a quality assurance review of the source test data and report. The review included the following activities:

- Review of the general text
- Check of calculation algorithms for correctness
- Review of CEMS data extraction
- Determination that all supporting documentation is present

A detailed manual review of the data was completed. Based on our data review, Shaw concludes the following:

- The source test procedures were followed, and all data were reduced and entered into the reporting spreadsheets properly.
- The instrument drift and other measures of instrument performance were deemed acceptable.

As a result, the data are considered to be acceptable.

## ***Limitations***

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The services described in this report were performed consistent with generally accepted professional consulting principles and practices. No other warranty, expressed or implied, is made. These services were performed consistent with our agreement with our client. This report is solely for the use and information of our client unless otherwise noted. Any reliance on this report by a third party is at such party's sole risk.

Opinions contained in this report apply to conditions existing when services were performed and are intended only for the client, purposes, locations, time frames, and project parameters indicated. We are not responsible for the impacts of any changes in environmental standards, practices, or regulations subsequent to performance of services. We do not warrant the accuracy of information supplied by others, or the use of segregated portions of this report.

***Appendix A***  
***Test Results***

## EMISSION RATE CALCULATION

Facility: Redwood Landfill  
Source : Flare A-50  
Test date: 6/7/2006

Temperature Std: 70  
Pressure Std: 29.92  
O2% correction: 15  
O2% correction: 3  
NOx & CO  
NMOC

| Time                           | 0811-0841<br>Run 1 |        | 0852-0922<br>Run 2 |        | 0931-1001<br>Run 3 |        | Avg   |        | Permit<br>Limits |
|--------------------------------|--------------------|--------|--------------------|--------|--------------------|--------|-------|--------|------------------|
|                                | Inlet              | Outlet | Inlet              | Outlet | Inlet              | Outlet | Inlet | Outlet |                  |
| Flare Temp (F)                 | 1,647              |        | 1,647              |        | 1,648              |        | 1,647 |        |                  |
| LFG rate                       | 874                |        | 872                |        | 870                |        | 872   |        |                  |
| Fuel factor                    | 9,147              |        | 9,147              |        | 9,147              |        | 9,147 |        |                  |
| Oxygen (%)                     |                    | 11.1   |                    | 11.0   |                    | 11.2   |       | 11.1   |                  |
| Flowrate (dscfm)               | 874                | 8,849  | 872                | 8,739  | 870                | 8,899  | 872   | 8,829  |                  |
| <b>Oxides of Nitrogen</b>      |                    |        |                    |        |                    |        |       |        |                  |
| MW = 46                        |                    |        |                    |        |                    |        |       |        |                  |
| NOx, ppm                       |                    | 18.7   |                    | 18.3   |                    | 18.1   |       | 18.4   |                  |
| NOx, ppm corrected to 15% O2   |                    | 11.3   |                    | 10.9   |                    | 11.0   |       | 11.1   | 15               |
| NOx, lb/hr                     |                    | 1.2    |                    | 1.1    |                    | 1.1    |       | 1.2    |                  |
| NOx, lb/MMBtu                  |                    | 0.043  |                    | 0.042  |                    | 0.042  |       | 0.043  | 0.06             |
| <b>Carbon Monoxide</b>         |                    |        |                    |        |                    |        |       |        |                  |
| MW = 28                        |                    |        |                    |        |                    |        |       |        |                  |
| CO, ppm                        |                    | 53.7   |                    | 53.6   |                    | 65.0   |       | 57.4   |                  |
| CO, ppm corrected to 15% O2    |                    | 32.3   |                    | 31.9   |                    | 39.5   |       | 34.6   | 123              |
| CO, lb/hr                      |                    | 2.1    |                    | 2.0    |                    | 2.5    |       | 2.2    |                  |
| CO, lb/MMBtu                   |                    | 0.076  |                    | 0.075  |                    | 0.093  |       | 0.081  | 0.30             |
| <b>Hydrocarbons as methane</b> |                    |        |                    |        |                    |        |       |        |                  |
| MW = 16                        |                    |        |                    |        |                    |        |       |        |                  |
| THC, ppm                       | NM                 | 1.1    | 512186             | 1.8    | NM                 | 3.7    |       | 2.2    |                  |
| THC, lb/hr                     |                    | 0.02   | 1,109              | 0.04   |                    | 0.08   |       | 0.05   |                  |
| Methane, ppm                   | NM                 | NM     | 511,730            | NM     | NM                 | NM     |       |        |                  |
| NMOC, ppm                      | NM                 | 1.1    | 456                | 1.8    | NM                 | 3.7    |       | 2.2    |                  |
| NMOC, ppm corrected to 3% O2   |                    | 2.0    |                    | 3.3    |                    | 6.8    |       | 4.0    | 30               |
| NMOC, lb/hr                    | NM                 | 0.02   | 1.0                | 0.04   | NM                 | 0.082  | 1.0   | 0.05   |                  |
| <b>Efficiency %</b>            |                    |        |                    |        |                    |        |       |        |                  |
| NMOC                           |                    | NM     |                    | 96.0%  |                    | NM     |       |        | 98               |

### Calculations

lb/hr = ppm x MW x Qs x 8.223x10<sup>-5</sup> / (Tstd + 460)

ppm corrected @ O2 = ppm measured x [(20.9 - O2% correction) / (20.9% O2 - O2 measured)]

lb/MMBtu = F-factor x MW x [1.3711x10<sup>-6</sup> / (Tstd + 460)] x [(20.9 / 20.9 - O2%)] x ppm

Efficiency (%) = in (lb/hr)-out (lb/hr)/in (lb/hr) x 100

NM = not measured

# Stack Gas Flow Rate ---- Fuel Usage

EPA Method 19

Facility Redwood Landfill

Source Flare A-50

Date 6/7/2006

|                               |            | Run 1 | Run 2 | Run 3 | Average |
|-------------------------------|------------|-------|-------|-------|---------|
| Gross Caloric Value (Btu/ft3) | Btu/ft3    | 519   | 519   | 519   | 519     |
| Stack Oxygen                  | %          | 11.1  | 11    | 11.2  | 11.1    |
| Fuel factor @ 68 F            | DSCF/MMBtu | 9,147 | 9,147 | 9,147 | 9,147   |

|                             |           |        |        |        |        |
|-----------------------------|-----------|--------|--------|--------|--------|
| LFG Rate (CFM) @ Tstd       | SCFM      | 874    | 872    | 870    | 872    |
| Fuel Flowrate (SCFH) @ Tstd | SCFH      | 52,440 | 52,320 | 52,200 | 52,320 |
| Million Btu per minute      | MMBtu/min | 0.454  | 0.453  | 0.452  | 0.453  |
| Heat Input (MMBtu/hr)       | MMBtu/hr  | 27.2   | 27.2   | 27.1   | 27.2   |
| Heat Input (MMBtu/day)      | MMBtu/day | 653.2  | 651.7  | 650.2  | 651.7  |

|                             |       |       |       |       |       |
|-----------------------------|-------|-------|-------|-------|-------|
| Stack Gas Flow Rate (dscfm) | dscfm | 8,849 | 8,739 | 8,899 | 8,829 |
|-----------------------------|-------|-------|-------|-------|-------|

## Calculations

$$\text{SCFM} = \text{CFM} * (\text{Tstd} + 460) * \text{gas PSIA} / 14.7 / (\text{gas F} + 460)$$

$$\text{SCFH} = \text{SCFM} * 60$$

$$\text{MMBtu/min} = \text{SCFM} * \text{Btu/ft}^3 / 1,000,000$$

$$\text{MMBtu/hr} = \text{MMBtu/min} * 60$$

$$\text{DSCFM} = \text{Fuel factor} * \text{MMBtu/min} * (20.9 / (20.9 - \text{O}_2\%)) * (\text{Tstd} + 460) / 528$$

## Notes

Corrected Fuel Rate (SCFM) @ Tstd data is from the system data aquisition

## EMISSION RATE CALCULATION

Facility: Redwood Landfill  
Source : Flare A-51  
Test date: 6/8/2006

Temperature Std: 70  
Pressure Std: 29.92  
O2% correction: 15  
O2% correction: 3  
NOx & CO  
NMOC

| Time                           | 0852-0922<br>Run 1 |        | 0933-1003<br>Run 2 |        | 1014-1144<br>Run 3 |        | Avg   |        | Permit |
|--------------------------------|--------------------|--------|--------------------|--------|--------------------|--------|-------|--------|--------|
|                                | Inlet              | Outlet | Inlet              | Outlet | Inlet              | Outlet | Inlet | Outlet | Limits |
| Flare Temp (F)                 | 1,448              |        | 1,452              |        | 1,451              |        | 1,450 |        |        |
| LFG rate                       | 2,618              |        | 2,601              |        | 2,591              |        | 2,603 |        |        |
| Fuel factor                    | 9,158              |        | 9,158              |        | 9,158              |        | 9,158 |        |        |
| Oxygen (%)                     |                    | 13.0   |                    | 12.9   |                    | 13.3   |       | 13.1   |        |
| Flowrate (dscfm)               | 2,618              | 31,397 | 2,601              | 30,804 | 2,591              | 32,300 | 2,603 | 31,500 |        |
| <b>Oxides of Nitrogen</b>      |                    |        |                    |        |                    |        |       |        |        |
| MW = 46                        |                    |        |                    |        |                    |        |       |        |        |
| NOx, ppm                       |                    | 17.0   |                    | 17.2   |                    | 15.7   |       | 16.6   |        |
| NOx, ppm corrected to 15% O2   |                    | 12.7   |                    | 12.7   |                    | 12.2   |       | 12.5   | 15     |
| NOx, lb/hr                     |                    | 3.8    |                    | 3.8    |                    | 3.6    |       | 3.7    |        |
| NOx, lb/MMBtu                  |                    | 0.049  |                    | 0.049  |                    | 0.047  |       | 0.048  | 0.06   |
| <b>Carbon Monoxide</b>         |                    |        |                    |        |                    |        |       |        |        |
| MW = 28                        |                    |        |                    |        |                    |        |       |        |        |
| CO, ppm                        |                    | 13.7   |                    | 11.6   |                    | 21.3   |       | 15.5   |        |
| CO, ppm corrected to 15% O2    |                    | 10.2   |                    | 8.6    |                    | 16.5   |       | 11.8   | 123    |
| CO, lb/hr                      |                    | 1.9    |                    | 1.6    |                    | 3.0    |       | 2.1    |        |
| CO, lb/MMBtu                   |                    | 0.024  |                    | 0.020  |                    | 0.039  |       | 0.028  | 0.30   |
| <b>Hydrocarbons as methane</b> |                    |        |                    |        |                    |        |       |        |        |
| MW = 16                        |                    |        |                    |        |                    |        |       |        |        |
| THC, ppm                       | NM                 | 2.3    | 488428             | 1.3    | NM                 | 5.0    |       | 2.9    |        |
| THC, lb/hr                     |                    | 0.18   | 3,154              | 0.10   |                    | 0.40   |       | 0.23   |        |
| Methane, ppm                   | NM                 | NM     | 488,215            | NM     | NM                 | NM     |       |        |        |
| NMOC, ppm                      | NM                 | 2.3    | 213                | 1.3    | NM                 | 5.0    | 213   | 2.9    |        |
| NMOC, ppm corrected to 3% O2   |                    | 5.2    |                    | 2.9    |                    | 11.8   |       | 6.6    | 30     |
| NMOC, lb/hr                    | NM                 | 0.18   | 1.4                | 0.10   | NM                 | 0.401  | 1.4   | 0.23   |        |
| <b>Efficiency %</b>            |                    |        |                    |        |                    |        |       |        |        |
| NMOC                           |                    | NM     |                    | 92.8%  |                    | NM     |       |        | 98     |

### Calculations

lb/hr = ppm x MW x Qs x 8.223x10<sup>-5</sup> / (Tstd + 460)

ppm corrected @ O2 = ppm measured x [(20.9 - O2% correction) / (20.9% O2 - O2 measured)]

lb/MMBtu = F-factor x MW x [1.3711x10<sup>-6</sup> / (Tstd + 460)] x [(20.9 / 20.9 - O2%)] x ppm

Efficiency (%) = in (lb/hr)-out (lb/hr)/in (lb/hr) x 100

NM = not measured

# Stack Gas Flow Rate ---- Fuel Usage

EPA Method 19

Facility Redwood Landfill

Source Flare A-51

Date 6/8/2006

|                                   |            | Run 1   | Run 2   | Run 3   | Average |
|-----------------------------------|------------|---------|---------|---------|---------|
| Gross Caloric Value (Btu/ft3)     | Btu/ft3    | 495     | 495     | 495     | 495     |
| Stack Oxygen                      | %          | 13      | 12.9    | 13.3    | 13.1    |
| Fuel factor @ 68 F                | DSCF/MMBtu | 9,158   | 9,158   | 9,158   | 9,158   |
| Corrected Fuel Rate (SCFM) @ Tstd | SCFM       | 2,618   | 2,601   | 2,591   | 2,603   |
| Fuel Flowrate (SCFH) @ Tstd       | SCFH       | 157,080 | 156,060 | 155,460 | 156,200 |
| Million Btu per minute            | MMBtu/min  | 1.296   | 1.287   | 1.283   | 1.289   |
| Heat Input (MMBtu/hr)             | MMBtu/hr   | 77.8    | 77.2    | 77.0    | 77.3    |
| Heat Input (MMBtu/day)            | MMBtu/day  | 1866.1  | 1854.0  | 1846.9  | 1855.7  |
| Stack Gas Flow Rate (dscfm)       | dscfm      | 31,397  | 30,804  | 32,300  | 31,500  |

## Calculations

$$\text{SCFM} = \text{CFM} * (\text{Tstd} + 460) * \text{gas PSIA} / 14.7 / (\text{gas F} + 460)$$

$$\text{SCFH} = \text{SCFM} * 60$$

$$\text{MMBtu/min} = \text{SCFM} * \text{Btu/ft}^3 / 1,000,000$$

$$\text{MMBtu/hr} = \text{MMBtu/min} * 60$$

$$\text{DSCFM} = \text{Fuel factor} * \text{MMBtu/min} * (20.9 / (20.9 - \text{O}_2\%)) * (\text{Tstd} + 460) / 528$$

## Notes

Corrected Fuel Rate (SCFM) @ Tstd data is from the system data acquisition

## ***Appendix B***

### ***Continuous Emission Monitoring***

- Strata<sup>TM</sup> Data Logger Printout
- Strip chart records
- Calibration Gas Certificates of Analysis

Run averages corrected for bias

Operator: B Johnston

Plant Name: Redwood Landfill

Location: Flare A50

| Run | O2<br>% | CO<br>ppm | THC<br>ppm | NOx<br>ppm |
|-----|---------|-----------|------------|------------|
| 1   | 11.1    | 53.7      | 1.1        | 18.7       |
| 2   | 11.0    | 53.6      | 1.8        | 18.3       |
| 3   | 11.2    | 65.1      | 3.7        | 18.1       |

Run averages corrected for bias

Operator: B Johnston

Plant Name: Redwood Landfill

Location: Flare A51

|     | O2   | CO   | THC | NOx  |
|-----|------|------|-----|------|
| Run | %    | ppm  | ppm | ppm  |
| 1   | 13.0 | 13.7 | 2.3 | 17.0 |
| 2   | 12.9 | 11.6 | 1.3 | 17.2 |
| 3   | 13.3 | 21.3 | 5.0 | 15.7 |

Operator: B Johnston  
Plant Name: Redwood Landfill  
Location: Flare A50  
Run Length: 30 minutes  
Sample Rate: 40 per minute  
Average Calibration Results: 1 minutes  
Automatic Sequence, Calibration Error: No  
Automatic Sequence, System Bias: No  
Max Response Time: Manual  
Max Response Time: 1 minutes  
Traverse During Run: No

| Active | Chan. | Analyte |       | Span  | Span  | Offset |
|--------|-------|---------|-------|-------|-------|--------|
|        |       | Name    | Units | Units | Volts | Volts  |
| Yes    | 1     | O2      | %     | 25    | 1     | 0      |
| Yes    | 3     | CO      | ppm   | 500   | .5    | 0      |
| Yes    | 4     | THC     | ppm   | 50    | 10    | 0      |
| Yes    | 5     | NOx     | ppm   | 50    | 10    | 0      |

Measurement System Preparation Table

| Gas Name | Reference Cylinder Numbers |   |   |   |   |   |   |   |   |    |    |    |    |    |    |    |
|----------|----------------------------|---|---|---|---|---|---|---|---|----|----|----|----|----|----|----|
|          | 1                          | 2 | 3 | 4 | 5 | 6 | 7 | 8 | 9 | 10 | 11 | 12 | 13 | 14 | 15 | 16 |
| O2       | Z                          | H |   |   |   |   |   |   |   |    |    |    |    |    |    |    |
| CO       | Z                          |   | H |   |   |   |   |   |   |    |    |    |    |    |    |    |
| THC      | Z                          |   |   | H |   |   |   |   |   |    |    |    |    |    |    |    |
| NOx      | Z                          |   |   |   | H |   |   |   |   |    |    |    |    |    |    |    |

Leak Check 0.1" at 14 in Vac. for 1 min.

Flare A50

~~Flare~~ R1 0811 - 0841

R2 0852 - 0922

R3 0931 - 0001

## STRATA Configuration Page 2

| Gas Name | Zero Reference Cylinder |      | Low Reference Cylinder |     |
|----------|-------------------------|------|------------------------|-----|
|          | No.                     | Conc | ID Number              |     |
| O2       | 1                       | 0    | 99                     | 999 |
| CO       | 1                       | 0    | 99                     | 999 |
| THC      | 1                       | 0    | 99                     | 999 |
| NOx      | 1                       | 0    | 99                     | 999 |

| Gas Name | Mid Reference Cylinder |      | High Reference Cylinder |  |
|----------|------------------------|------|-------------------------|--|
|          | No.                    | Conc | ID Number               |  |
| O2       | 99                     | 999  | 999999                  |  |
| CO       | 99                     | 999  |                         |  |
| THC      | 99                     | 999  |                         |  |
| NOx      | 99                     | 999  |                         |  |

| Seq Num | Calibration Error Test Sequence |    |     |     |
|---------|---------------------------------|----|-----|-----|
|         | O2                              | CO | THC | NOx |

1  
2  
3  
4  
5  
6  
7  
8  
9  
10  
11  
12  
13  
14  
15

| Seq Num | Calibration Error Valve Sequence |   |   |   |   |   |   |   |   |    |    |    |    |    |    |    |
|---------|----------------------------------|---|---|---|---|---|---|---|---|----|----|----|----|----|----|----|
|         | 1                                | 2 | 3 | 4 | 5 | 6 | 7 | 8 | 9 | 10 | 11 | 12 | 13 | 14 | 15 | 16 |

1  
2  
3  
4  
5  
6  
7  
8  
9  
10  
11  
12  
13  
14  
15

# STRATA Configuration Page 3

| Seq<br>Num | System Bias Check Sequence |      |      |      |
|------------|----------------------------|------|------|------|
|            | O2                         | CO   | THC  | NOx  |
| 1          | Zero                       | Zero | Zero | Zero |
| 2          | High                       | High | High | High |
| 3          |                            |      |      |      |
| 4          |                            |      |      |      |
| 5          |                            |      |      |      |
| 6          |                            |      |      |      |
| 7          |                            |      |      |      |
| 8          |                            |      |      |      |
| 9          |                            |      |      |      |
| 10         |                            |      |      |      |
| 11         |                            |      |      |      |
| 12         |                            |      |      |      |
| 13         |                            |      |      |      |
| 14         |                            |      |      |      |
| 15         |                            |      |      |      |

| Seq<br>Num | System Bias Valve Sequence |   |   |   |   |   |   |   |   |    |    |    |    |    |    |    |
|------------|----------------------------|---|---|---|---|---|---|---|---|----|----|----|----|----|----|----|
|            | 1                          | 2 | 3 | 4 | 5 | 6 | 7 | 8 | 9 | 10 | 11 | 12 | 13 | 14 | 15 | 16 |
| 1          | X                          |   |   |   |   |   |   |   |   |    |    |    |    |    |    |    |
| 2          |                            | X | X | X | X |   |   |   |   |    |    |    |    |    |    |    |
| 3          |                            |   |   |   |   |   |   |   |   |    |    |    |    |    |    |    |
| 4          |                            |   |   |   |   |   |   |   |   |    |    |    |    |    |    |    |
| 5          |                            |   |   |   |   |   |   |   |   |    |    |    |    |    |    |    |
| 6          |                            |   |   |   |   |   |   |   |   |    |    |    |    |    |    |    |
| 7          |                            |   |   |   |   |   |   |   |   |    |    |    |    |    |    |    |
| 8          |                            |   |   |   |   |   |   |   |   |    |    |    |    |    |    |    |
| 9          |                            |   |   |   |   |   |   |   |   |    |    |    |    |    |    |    |
| 10         |                            |   |   |   |   |   |   |   |   |    |    |    |    |    |    |    |
| 11         |                            |   |   |   |   |   |   |   |   |    |    |    |    |    |    |    |
| 12         |                            |   |   |   |   |   |   |   |   |    |    |    |    |    |    |    |
| 13         |                            |   |   |   |   |   |   |   |   |    |    |    |    |    |    |    |
| 14         |                            |   |   |   |   |   |   |   |   |    |    |    |    |    |    |    |
| 15         |                            |   |   |   |   |   |   |   |   |    |    |    |    |    |    |    |

STRATA Configuration End

## Initial System Bias Check, Run 1 STRATA Version 1.1

|                     | O2<br>% | CO<br>ppm | THC<br>ppm | NOx<br>ppm |
|---------------------|---------|-----------|------------|------------|
| 06-07-2006 07:58:33 | 3.773   | -1.43     | -0.034     | 2.647      |
| 06-07-2006 07:59:34 | 0.154   | -2.23     | -0.080     | 0.010      |
| 06-07-2006 08:00:33 | 0.107   | -0.90     | -0.079     | 0.011      |
| 06-07-2006 08:01:33 | 0.209   | -0.96     | 3.201      | 3.456      |
| 06-07-2006 08:02:34 | 17.051  | 84.31     | 44.708     | 43.742     |
| 06-07-2006 08:03:33 | 19.303  | 90.67     | 44.767     | 45.188     |
| 06-07-2006 08:04:34 | 19.353  | 90.16     | 44.736     | 45.384     |
| 06-07-2006 08:05:33 | 19.368  | 90.38     | 44.751     | 45.385     |
| 06-07-2006 08:06:33 | 19.376  | 90.12     | 44.772     | 45.041     |

## Initial System Bias Check for Run 1

Operator: B Johnston  
Plant Name: Redwood Landfill  
Location: Flare A50

## Reference Cylinder Numbers

|     | Zero | Span     |
|-----|------|----------|
| O2  |      | CC8445   |
| CO  |      | AAL19769 |
| THC |      | AAL4576  |
| NOx |      | AAL19769 |

| Date/Time             | 06-07-2006 | 08:07:28 | PASSED |
|-----------------------|------------|----------|--------|
| Analyte               | O2         | CO       | THC    |
| Units                 | %          | ppm      | ppm    |
| Zero Ref Cyl          | 0.000      | 0.00     | 0.000  |
| Zero Cal              | 0.000      | 0.00     | 0.000  |
| Zero Avg              | 0.098      | -1.03    | -0.078 |
| Zero Bias%            | 0.4%       | 0.2%     | 0.2%   |
| Zero Drift%           |            |          |        |
| Span Ref Cyl          | 19.040     | 90.10    | 45.000 |
| Span Cal              | 19.040     | 90.10    | 45.000 |
| Span Avg              | 19.381     | 90.14    | 44.771 |
| Span Bias%            | 1.4%       | 0.0%     | 0.5%   |
| Span Drift%           |            |          |        |
| System Bias Check End |            |          |        |

Test Run 1 STRATA Version 1.1

|                                | O2<br>% | CO<br>ppm | THC<br>ppm | NOx<br>ppm |
|--------------------------------|---------|-----------|------------|------------|
| 06-07-2006 08:08:28            | 14.557  | 37.14     | 6.647      | 23.338     |
| 06-07-2006 08:09:29            | 11.473  | 37.27     | 0.076      | 18.787     |
| 06-07-2006 08:10:28            | 11.974  | 82.16     | 3.118      | 17.129     |
| 06-07-2006 08:11:29            | 11.328  | 17.85     | -0.074     | 18.865     |
| Begin calculating run averages |         |           |            |            |
| 06-07-2006 08:12:28            | 11.920  | 32.25     | -0.074     | 17.515     |
| 06-07-2006 08:13:28            | 11.426  | 24.90     | -0.075     | 18.824     |
| 06-07-2006 08:14:29            | 11.131  | 56.38     | 0.120      | 18.752     |
| 06-07-2006 08:15:28            | 11.811  | 64.21     | 1.022      | 17.967     |
| 06-07-2006 08:16:29            | 10.982  | 38.32     | 0.191      | 19.522     |
| 06-07-2006 08:17:29            | 11.804  | 52.62     | 0.062      | 17.137     |
| 06-07-2006 08:18:28            | 11.871  | 89.55     | 3.706      | 18.284     |
| 06-07-2006 08:19:29            | 11.222  | 54.57     | 0.497      | 18.554     |
| 06-07-2006 08:20:28            | 11.975  | 111.67    | 6.487      | 17.582     |
| 06-07-2006 08:21:28            | 11.516  | 31.92     | -0.074     | 18.595     |
| 06-07-2006 08:22:29            | 10.570  | 107.73    | 2.069      | 20.223     |
| 06-07-2006 08:23:28            | 11.071  | 38.45     | 0.072      | 18.446     |
| 06-07-2006 08:24:29            | 11.790  | 80.33     | 2.306      | 18.317     |
| 06-07-2006 08:25:28            | 11.467  | 57.66     | 0.218      | 18.565     |
| 06-07-2006 08:26:28            | 10.511  | 47.48     | 0.104      | 20.485     |
| 06-07-2006 08:27:29            | 11.369  | 66.41     | 0.612      | 17.991     |
| 06-07-2006 08:28:28            | 11.853  | 23.79     | -0.075     | 17.585     |
| 06-07-2006 08:29:28            | 11.812  | 79.80     | 1.726      | 18.355     |
| 06-07-2006 08:30:29            | 10.899  | 34.04     | -0.054     | 19.537     |
| 06-07-2006 08:31:28            | 11.046  | 79.48     | 2.023      | 18.948     |
| 06-07-2006 08:32:29            | 11.865  | 71.86     | 1.877      | 17.651     |
| 06-07-2006 08:33:29            | 11.142  | 24.88     | -0.009     | 19.065     |
| 06-07-2006 08:34:28            | 11.469  | 21.96     | -0.075     | 18.145     |
| 06-07-2006 08:35:29            | 10.726  | 18.53     | -0.075     | 20.214     |
| 06-07-2006 08:36:28            | 10.387  | 2.38      | -0.075     | 20.460     |
| 06-07-2006 08:37:29            | 10.543  | 79.21     | 2.402      | 20.357     |
| 06-07-2006 08:38:29            | 10.544  | 48.57     | 1.101      | 20.066     |
| 06-07-2006 08:39:28            | 11.300  | 57.31     | 0.601      | 18.429     |
| 06-07-2006 08:40:29            | 11.147  | 24.58     | -0.075     | 18.476     |
| 06-07-2006 08:41:28            | 11.566  | 35.76     | 0.036      | 18.285     |
| Run Averages                   |         |           |            |            |
|                                | O2<br>% | CO<br>ppm | THC<br>ppm | NOx<br>ppm |
| 06-07-2006 08:41:53            | 11.268  | 53.93     | 0.989      | 18.799     |
| Operator: B Johnston           |         |           |            |            |
| Plant Name: Redwood Landfill   |         |           |            |            |
| Location: Flare A50            |         |           |            |            |
| Test Run 1 End                 |         |           |            |            |

# Final System Bias Check, Run 1 STRATA Version 1.1

|                     | O2<br>% | CO<br>ppm | THC<br>ppm | NOx<br>ppm |
|---------------------|---------|-----------|------------|------------|
| 06-07-2006 08:43:15 | 2.627   | -2.30     | -0.076     | 0.018      |
| 06-07-2006 08:44:16 | 0.172   | -4.97     | -0.076     | 0.015      |
| 06-07-2006 08:45:15 | 0.133   | 1.38      | -0.075     | 0.015      |
| 06-07-2006 08:46:15 | 3.116   | 23.88     | 13.577     | 14.900     |
| 06-07-2006 08:47:14 | 18.601  | 93.31     | 45.558     | 43.832     |
| 06-07-2006 08:48:15 | 19.254  | 90.48     | 45.004     | 44.997     |
| 06-07-2006 08:49:16 | 19.284  | 90.59     | 45.114     | 45.088     |

## Final System Bias Check for Run 1

Operator: B Johnston  
Plant Name: Redwood Landfill  
Location: Flare A50

### Reference Cylinder Numbers

|     | Zero | Span     |
|-----|------|----------|
| O2  |      | CC8445   |
| CO  |      | AAL19769 |
| THC |      | AAL4576  |
| NOx |      | AAL19769 |

| Date/Time             | 06-07-2006 | 08:49:22 | PASSED |
|-----------------------|------------|----------|--------|
| Analyte               | O2         | CO       | THC    |
| Units                 | %          | ppm      | ppm    |
| Zero Ref Cyl          | 0.000      | 0.00     | 0.000  |
| Zero Cal              | 0.000      | 0.00     | 0.000  |
| Zero Avg              | 0.125      | 1.36     | -0.075 |
| Zero Bias%            | 0.5%       | 0.3%     | 0.2%   |
| Zero Drift%           | 0.1%       | 0.5%     | 0.0%   |
| Span Ref Cyl          | 19.040     | 90.10    | 45.000 |
| Span Cal              | 19.040     | 90.10    | 45.000 |
| Span Avg              | 19.284     | 90.60    | 45.111 |
| Span Bias%            | 1.0%       | 0.1%     | 0.2%   |
| Span Drift%           | -0.4%      | 0.1%     | 0.7%   |
| Ini Zero Avg          | 0.098      | -1.03    | -0.078 |
| Ini Span Avg          | 19.381     | 90.14    | 44.771 |
| Run Avg               | 11.268     | 53.93    | 0.989  |
| Co                    | 0.112      | 0.16     | -0.077 |
| Cm                    | 19.333     | 90.37    | 44.941 |
| Correct Avg           | 11.052     | 53.71    | 1.066  |
| System Bias Check End |            |          |        |

Test Run 2 STRATA Version 1.1

|                                | O2<br>% | CO<br>ppm | THC<br>ppm | NOx<br>ppm |
|--------------------------------|---------|-----------|------------|------------|
| 06-07-2006 08:50:23            | 13.814  | 35.95     | 5.200      | 22.229     |
| 06-07-2006 08:51:24            | 10.870  | 67.85     | 3.742      | 19.342     |
| Begin calculating run averages |         |           |            |            |
| 06-07-2006 08:52:23            | 11.117  | 17.83     | 0.116      | 17.661     |
| 06-07-2006 08:53:24            | 11.047  | 25.44     | 0.440      | 19.358     |
| 06-07-2006 08:54:23            | 10.364  | 109.68    | 4.543      | 19.926     |
| 06-07-2006 08:55:23            | 10.922  | 62.07     | 1.227      | 19.235     |
| 06-07-2006 08:56:24            | 10.652  | 28.39     | 1.592      | 19.606     |
| 06-07-2006 08:57:23            | 11.168  | 14.17     | -0.031     | 18.446     |
| 06-07-2006 08:58:23            | 11.256  | 42.20     | 0.705      | 18.101     |
| 06-07-2006 08:59:24            | 10.935  | 122.95    | 5.206      | 18.795     |
| 06-07-2006 09:00:23            | 11.372  | 54.61     | 1.669      | 18.002     |
| 06-07-2006 09:01:24            | 11.615  | 54.32     | 0.939      | 17.607     |
| 06-07-2006 09:02:23            | 10.906  | 72.38     | 1.030      | 18.917     |
| 06-07-2006 09:03:23            | 11.131  | 41.62     | 0.219      | 18.509     |
| 06-07-2006 09:04:24            | 10.845  | 13.74     | 0.001      | 19.053     |
| 06-07-2006 09:05:23            | 11.381  | 27.14     | 0.015      | 17.851     |
| 06-07-2006 09:06:23            | 11.186  | 36.20     | 0.107      | 18.365     |
| 06-07-2006 09:07:22            | 11.421  | 30.66     | -0.034     | 17.667     |
| 06-07-2006 09:08:23            | 11.505  | 25.27     | -0.071     | 18.065     |
| 06-07-2006 09:09:24            | 11.260  | 50.28     | 0.566      | 18.343     |
| 06-07-2006 09:10:23            | 10.601  | 79.16     | 2.195      | 19.684     |
| 06-07-2006 09:11:23            | 10.374  | 33.36     | 1.147      | 19.797     |
| 06-07-2006 09:12:24            | 11.861  | 111.29    | 9.425      | 16.580     |
| 06-07-2006 09:13:23            | 11.408  | 33.52     | 0.337      | 17.929     |
| 06-07-2006 09:14:23            | 11.497  | 41.73     | 0.665      | 18.342     |
| 06-07-2006 09:15:24            | 10.654  | 31.35     | 0.558      | 19.545     |
| 06-07-2006 09:16:23            | 11.350  | 41.92     | 0.577      | 17.444     |
| 06-07-2006 09:17:24            | 11.812  | 58.78     | 0.552      | 17.032     |
| 06-07-2006 09:18:23            | 12.206  | 129.32    | 7.916      | 15.881     |
| 06-07-2006 09:19:23            | 11.030  | 43.89     | 0.850      | 19.311     |
| 06-07-2006 09:20:24            | 11.429  | 75.16     | 2.262      | 17.533     |
| 06-07-2006 09:21:23            | 11.867  | 112.13    | 7.106      | 17.546     |
| Run Averages                   | O2<br>% | CO<br>ppm | THC<br>ppm | NOx<br>ppm |
| 06-07-2006 09:22:05            | 11.206  | 54.37     | 1.728      | 18.357     |

Operator: B Johnston  
Plant Name: Redwood Landfill  
Location: Flare A50  
Test Run 2 End

Final System Bias Check, Run 2 STRATA Version 1.1

|            |          | O2<br>% | CO<br>ppm | THC<br>ppm | NOx<br>ppm |
|------------|----------|---------|-----------|------------|------------|
| 06-07-2006 | 09:23:26 | 2.844   | 2.83      | -0.067     | 0.014      |
| 06-07-2006 | 09:24:25 | 0.154   | -0.37     | -0.081     | 0.011      |
| 06-07-2006 | 09:25:26 | 0.104   | -0.69     | -0.081     | 0.011      |
| 06-07-2006 | 09:26:27 | 13.243  | 75.75     | 39.187     | 39.924     |
| 06-07-2006 | 09:27:26 | 19.092  | 91.48     | 44.995     | 44.734     |
| 06-07-2006 | 09:28:27 | 19.185  | 91.60     | 44.963     | 44.722     |

Final System Bias Check for Run 2

Operator: B Johnston  
 Plant Name: Redwood Landfill  
 Location: Flare A50

Reference Cylinder Numbers

|     | Zero | Span     |
|-----|------|----------|
| O2  |      | CC8445   |
| CO  |      | AAL19769 |
| THC |      | AAL4576  |
| NOx |      | AAL19769 |

| Date/Time             | 06-07-2006 |       | 09:28:27 |        | PASSED |
|-----------------------|------------|-------|----------|--------|--------|
| Analyte               | O2         | CO    | THC      | NOx    |        |
| Units                 | %          | ppm   | ppm      | ppm    |        |
| Zero Ref Cyl          | 0.000      | 0.00  | 0.000    | 0.000  |        |
| Zero Cal              | 0.000      | 0.00  | 0.000    | 0.000  |        |
| Zero Avg              | 0.105      | -0.70 | -0.081   | 0.011  |        |
| Zero Bias%            | 0.4%       | 0.1%  | 0.2%     | 0.0%   |        |
| Zero Drift%           | -0.1%      | -0.4% | 0.0%     | 0.0%   |        |
| Span Ref Cyl          | 19.040     | 90.10 | 45.000   | 44.800 |        |
| Span Cal              | 19.040     | 90.10 | 45.000   | 44.800 |        |
| Span Avg              | 19.182     | 91.64 | 44.969   | 44.724 |        |
| Span Bias%            | 0.6%       | 0.3%  | 0.1%     | 0.2%   |        |
| Span Drift%           | -0.4%      | 0.2%  | -0.3%    | -0.7%  |        |
| Ini Zero Avg          | 0.125      | 1.36  | -0.075   | 0.015  |        |
| Ini Span Avg          | 19.284     | 90.60 | 45.111   | 45.091 |        |
| Run Avg               | 11.206     | 54.37 | 1.728    | 18.357 |        |
| Co                    | 0.115      | 0.33  | -0.078   | 0.013  |        |
| Cm                    | 19.233     | 91.12 | 45.040   | 44.907 |        |
| Correct Avg           | 11.046     | 53.63 | 1.802    | 18.305 |        |
| System Bias Check End |            |       |          |        |        |

Test Run 3 STRATA Version 1.1

|                                |  | O2<br>% | CO<br>ppm | THC<br>ppm | NOx<br>ppm |
|--------------------------------|--|---------|-----------|------------|------------|
| 06-07-2006 09:29:28            |  | 13.085  | 43.97     | 6.085      | 21.992     |
| 06-07-2006 09:30:28            |  | 11.089  | 29.28     | 0.109      | 18.323     |
| 06-07-2006 09:31:27            |  | 11.249  | 43.48     | 1.176      | 18.853     |
| Begin calculating run averages |  |         |           |            |            |
| 06-07-2006 09:32:28            |  | 10.632  | 80.97     | 6.031      | 19.212     |
| 06-07-2006 09:33:27            |  | 10.730  | 56.19     | 1.086      | 18.959     |
| 06-07-2006 09:34:28            |  | 11.276  | 49.26     | 0.289      | 18.065     |
| 06-07-2006 09:35:28            |  | 11.264  | 56.52     | 1.548      | 18.508     |
| 06-07-2006 09:36:27            |  | 11.507  | 98.61     | 4.786      | 17.367     |
| 06-07-2006 09:37:28            |  | 11.576  | 44.93     | 1.206      | 17.898     |
| 06-07-2006 09:38:28            |  | 10.895  | 24.13     | 0.291      | 18.889     |
| 06-07-2006 09:39:27            |  | 12.071  | 141.85    | 8.607      | 16.447     |
| 06-07-2006 09:40:28            |  | 10.806  | 34.77     | 2.044      | 19.092     |
| 06-07-2006 09:41:27            |  | 10.601  | 41.87     | 1.537      | 19.935     |
| 06-07-2006 09:42:28            |  | 10.578  | 29.53     | 0.352      | 18.737     |
| 06-07-2006 09:43:28            |  | 11.002  | 52.78     | 1.676      | 19.242     |
| 06-07-2006 09:44:27            |  | 10.696  | 21.75     | 0.591      | 19.278     |
| 06-07-2006 09:45:28            |  | 10.811  | 46.56     | 1.825      | 19.141     |
| 06-07-2006 09:46:28            |  | 10.698  | 18.20     | 0.530      | 19.340     |
| 06-07-2006 09:47:27            |  | 10.593  | 53.85     | 1.077      | 19.488     |
| 06-07-2006 09:48:28            |  | 10.834  | 52.40     | 1.092      | 18.406     |
| 06-07-2006 09:49:27            |  | 10.963  | 38.63     | 1.183      | 18.569     |
| 06-07-2006 09:50:28            |  | 11.108  | 15.34     | 0.285      | 18.674     |
| 06-07-2006 09:51:28            |  | 11.328  | 70.48     | 1.724      | 17.379     |
| 06-07-2006 09:52:27            |  | 11.613  | 61.89     | 3.560      | 18.180     |
| 06-07-2006 09:53:28            |  | 11.467  | 54.68     | 1.094      | 17.669     |
| 06-07-2006 09:54:28            |  | 12.500  | 154.25    | 18.466     | 16.057     |
| 06-07-2006 09:55:27            |  | 12.409  | 142.76    | 12.391     | 16.334     |
| 06-07-2006 09:56:28            |  | 11.763  | 101.21    | 5.461      | 17.434     |
| 06-07-2006 09:57:27            |  | 11.836  | 39.09     | 0.253      | 17.028     |
| 06-07-2006 09:58:28            |  | 11.867  | 109.05    | 6.140      | 17.614     |
| 06-07-2006 09:59:28            |  | 12.132  | 139.24    | 13.128     | 16.607     |
| 06-07-2006 10:00:27            |  | 12.189  | 114.16    | 9.299      | 16.352     |
| 06-07-2006 10:01:28            |  | 11.574  | 43.39     | 1.563      | 17.423     |
| Run Averages                   |  | O2<br>% | CO<br>ppm | THC<br>ppm | NOx<br>ppm |
| 06-07-2006 10:01:31            |  | 11.313  | 66.33     | 3.628      | 18.106     |

Operator: B Johnston  
Plant Name: Redwood Landfill  
Location: Flare A50  
Test Run 3 End

# Final System Bias Check, Run 3 STRATA Version 1.1

|                     | O2<br>% | CO<br>ppm | THC<br>ppm | NOx<br>ppm |
|---------------------|---------|-----------|------------|------------|
| 06-07-2006 10:02:40 | 2.907   | 4.06      | 0.268      | 0.008      |
| 06-07-2006 10:03:39 | 0.129   | 0.13      | -0.087     | 0.006      |
| 06-07-2006 10:04:40 | 0.080   | 0.13      | -0.086     | 0.006      |
| 06-07-2006 10:05:40 | 8.934   | 44.15     | 29.831     | 29.022     |
| 06-07-2006 10:06:39 | 18.962  | 92.55     | 44.922     | 44.554     |
| 06-07-2006 10:07:40 | 19.130  | 92.32     | 45.006     | 44.824     |

## Final System Bias Check for Run 3

Operator: B Johnston  
 Plant Name: Redwood Landfill  
 Location: Flare A50

### Reference Cylinder Numbers

|     | Zero | Span     |
|-----|------|----------|
| O2  |      | CC8445   |
| CO  |      | AAL19769 |
| THC |      | AAL4576  |
| NOx |      | AAL19769 |

| Date/Time             | 06-07-2006 | 10:08:17 | PASSED |
|-----------------------|------------|----------|--------|
| Analyte               | O2         | CO       | THC    |
| Units                 | %          | ppm      | ppm    |
| Zero Ref Cyl          | 0.000      | 0.00     | 0.000  |
| Zero Cal              | 0.000      | 0.00     | 0.000  |
| Zero Avg              | 0.077      | 0.16     | -0.086 |
| Zero Bias%            | 0.3%       | 0.0%     | 0.2%   |
| Zero Drift%           | -0.1%      | 0.2%     | 0.0%   |
| Span Ref Cyl          | 19.040     | 90.10    | 45.000 |
| Span Cal              | 19.040     | 90.10    | 45.000 |
| Span Avg              | 19.143     | 92.31    | 45.053 |
| Span Bias%            | 0.4%       | 0.4%     | 0.1%   |
| Span Drift%           | -0.2%      | 0.1%     | 0.2%   |
| Ini Zero Avg          | 0.105      | -0.70    | -0.081 |
| Ini Span Avg          | 19.182     | 91.64    | 44.969 |
| Run Avg               | 11.313     | 66.33    | 3.628  |
| Co                    | 0.091      | -0.27    | -0.084 |
| Cm                    | 19.163     | 91.97    | 45.011 |
| Correct Avg           | 11.203     | 65.05    | 3.704  |
| System Bias Check End |            |          |        |

STRATA Configuration Page 1  
06-08-2006 08:15:47  
File Name: C:\STRATA\REDWOOD.STR

Operator: B Johnston  
Plant Name: Redwood Landfill  
Location: Flare A51  
Run Length: 30 minutes  
Sample Rate: 40 per minute  
Average Calibration Results: 1 minutes  
Automatic Sequence, Calibration Error: No  
Automatic Sequence, System Bias: No  
Max Response Time: Manual  
Max Response Time: 1 minutes  
Traverse During Run: No

| Active | Chan. | Analyte Name | Units | Span Units | Span Volts | Offset Volts |
|--------|-------|--------------|-------|------------|------------|--------------|
| Yes    | 1     | O2           | %     | 25         | 1          | 0            |
| Yes    | 3     | CO           | ppm   | 100        | .5         | 0            |
| Yes    | 4     | THC          | ppm   | 50         | 10         | 0            |
| Yes    | 5     | NOx          | ppm   | 50         | 10         | 0            |

Measurement System Preparation Table

| Gas Name | 1 | 2 | 3 | 4 | 5 | 6 | 7 | 8 | 9 | 10 | 11 | 12 | 13 | 14 | 15 | 16 |
|----------|---|---|---|---|---|---|---|---|---|----|----|----|----|----|----|----|
| O2       | Z | H |   |   |   |   |   |   |   |    |    |    |    |    |    |    |
| CO       | Z |   | H |   |   |   |   |   |   |    |    |    |    |    |    |    |
| THC      | Z |   |   | H |   |   |   |   |   |    |    |    |    |    |    |    |
| NOx      | Z |   |   |   | H |   |   |   |   |    |    |    |    |    |    |    |

Leak Check 0.1" at 14" vacuum for 1 min

R1 0852 - 0922  
R2 0933 - 1003  
R3 1014 - 1044

## STRATA Configuration Page 2

| Gas Name | Zero Reference Cylinder |      | Low Reference Cylinder |  |
|----------|-------------------------|------|------------------------|--|
|          | No.                     | Conc | ID Number              |  |
| O2       | 1                       | 0    |                        |  |
| CO       | 1                       | 0    |                        |  |
| THC      | 1                       | 0    |                        |  |
| NOx      | 1                       | 0    |                        |  |

| Gas Name | Mid Reference Cylinder |      | High Reference Cylinder |  |
|----------|------------------------|------|-------------------------|--|
|          | No.                    | Conc | ID Number               |  |
| O2       | 99                     | 999  | 999999                  |  |
| CO       | 99                     | 999  |                         |  |
| THC      | 99                     | 999  |                         |  |
| NOx      | 99                     | 999  |                         |  |

| Seq Num | Calibration Error Test Sequence |    |     |     |
|---------|---------------------------------|----|-----|-----|
|         | O2                              | CO | THC | NOx |

1  
2  
3  
4  
5  
6  
7  
8  
9  
10  
11  
12  
13  
14  
15

| Seq Num | Calibration Error Valve Sequence |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
|---------|----------------------------------|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|
|---------|----------------------------------|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|

|  | 1 | 2 | 3 | 4 | 5 | 6 | 7 | 8 | 9 | 10 | 11 | 12 | 13 | 14 | 15 | 16 |
|--|---|---|---|---|---|---|---|---|---|----|----|----|----|----|----|----|
|--|---|---|---|---|---|---|---|---|---|----|----|----|----|----|----|----|

1  
2  
3  
4  
5  
6  
7  
8  
9  
10  
11  
12  
13  
14  
15

# STRATA Configuration Page 3

| Seq<br>Num | System Bias Check Sequence |      |      |      |
|------------|----------------------------|------|------|------|
|            | O2                         | CO   | THC  | NOx  |
| 1          | Zero                       | Zero | Zero | Zero |
| 2          | High                       | High | High | High |
| 3          |                            |      |      |      |
| 4          |                            |      |      |      |
| 5          |                            |      |      |      |
| 6          |                            |      |      |      |
| 7          |                            |      |      |      |
| 8          |                            |      |      |      |
| 9          |                            |      |      |      |
| 10         |                            |      |      |      |
| 11         |                            |      |      |      |
| 12         |                            |      |      |      |
| 13         |                            |      |      |      |
| 14         |                            |      |      |      |
| 15         |                            |      |      |      |

| Seq<br>Num | System Bias Valve Sequence |   |   |   |   |   |   |   |   |    |    |    |    |    |    |    |
|------------|----------------------------|---|---|---|---|---|---|---|---|----|----|----|----|----|----|----|
|            | 1                          | 2 | 3 | 4 | 5 | 6 | 7 | 8 | 9 | 10 | 11 | 12 | 13 | 14 | 15 | 16 |
| 1          | X                          |   |   |   |   |   |   |   |   |    |    |    |    |    |    |    |
| 2          |                            | X | X | X | X |   |   |   |   |    |    |    |    |    |    |    |
| 3          |                            |   |   |   |   |   |   |   |   |    |    |    |    |    |    |    |
| 4          |                            |   |   |   |   |   |   |   |   |    |    |    |    |    |    |    |
| 5          |                            |   |   |   |   |   |   |   |   |    |    |    |    |    |    |    |
| 6          |                            |   |   |   |   |   |   |   |   |    |    |    |    |    |    |    |
| 7          |                            |   |   |   |   |   |   |   |   |    |    |    |    |    |    |    |
| 8          |                            |   |   |   |   |   |   |   |   |    |    |    |    |    |    |    |
| 9          |                            |   |   |   |   |   |   |   |   |    |    |    |    |    |    |    |
| 10         |                            |   |   |   |   |   |   |   |   |    |    |    |    |    |    |    |
| 11         |                            |   |   |   |   |   |   |   |   |    |    |    |    |    |    |    |
| 12         |                            |   |   |   |   |   |   |   |   |    |    |    |    |    |    |    |
| 13         |                            |   |   |   |   |   |   |   |   |    |    |    |    |    |    |    |
| 14         |                            |   |   |   |   |   |   |   |   |    |    |    |    |    |    |    |
| 15         |                            |   |   |   |   |   |   |   |   |    |    |    |    |    |    |    |

STRATA Configuration End

## Initial System Bias Check, Run 1 STRATA Version 1.1

|                     | O2<br>% | CO<br>ppm | THC<br>ppm | NOx<br>ppm |
|---------------------|---------|-----------|------------|------------|
| 06-08-2006 08:24:39 | 4.564   | 2.49      | 1.456      | 1.777      |
| 06-08-2006 08:25:40 | 0.203   | 0.06      | -0.076     | 0.010      |
| 06-08-2006 08:26:39 | 0.126   | -0.10     | -0.076     | 0.010      |
| 06-08-2006 08:27:39 | 0.108   | -0.12     | -0.076     | 0.010      |
| 06-08-2006 08:28:38 | 5.376   | 16.49     | 22.318     | 21.111     |
| 06-08-2006 08:29:39 | 18.672  | 90.35     | 44.994     | 44.343     |
| 06-08-2006 08:30:39 | 19.046  | 90.49     | 45.015     | 45.420     |
| 06-08-2006 08:31:38 | 19.077  | 90.73     | 45.092     | 44.959     |
| 06-08-2006 08:32:39 | 19.086  | 90.21     | 45.127     | 44.844     |

## Initial System Bias Check for Run 1

Operator: B Johnston  
 Plant Name: Redwood Landfill  
 Location: Flare A51

## Reference Cylinder Numbers

|     | Zero | Span     |
|-----|------|----------|
| O2  |      | CC8445   |
| CO  |      | AAL19769 |
| THC |      | AAL4576  |
| NOx |      | AAL19769 |

| Date/Time             | 06-08-2006 | 08:33:24 | PASSED |
|-----------------------|------------|----------|--------|
| Analyte               | O2         | CO       | THC    |
| Units                 | %          | ppm      | ppm    |
| Zero Ref Cyl          | 0.000      | 0.00     | 0.000  |
| Zero Cal              | 0.000      | 0.00     | 0.000  |
| Zero Avg              | 0.105      | -0.10    | -0.076 |
| Zero Bias%            | 0.4%       | 0.1%     | 0.2%   |
| Zero Drift%           |            |          |        |
| Span Ref Cyl          | 19.040     | 90.10    | 45.000 |
| Span Cal              | 19.040     | 90.10    | 45.000 |
| Span Avg              | 19.088     | 90.46    | 45.090 |
| Span Bias%            | 0.2%       | 0.4%     | 0.2%   |
| Span Drift%           |            |          |        |
| System Bias Check End |            |          |        |

Test Run 1 STRATA Version 1.1

|                                |          | O2<br>% | CO<br>ppm | THC<br>ppm | NOx<br>ppm |
|--------------------------------|----------|---------|-----------|------------|------------|
| 06-08-2006                     | 08:34:25 | 15.210  | 37.68     | 11.697     | 18.309     |
| 06-08-2006                     | 08:35:25 | 13.938  | 27.99     | 9.898      | 14.770     |
| 06-08-2006                     | 08:36:24 | 15.743  | 15.79     | 27.653     | 7.678      |
| 06-08-2006                     | 08:37:25 | 20.362  | 21.90     | 49.986     | 1.113      |
| 06-08-2006                     | 08:38:24 | 13.266  | 34.32     | 43.522     | 19.376     |
| 06-08-2006                     | 08:39:25 | 12.639  | 12.01     | 15.781     | 17.965     |
| 06-08-2006                     | 08:40:24 | 13.237  | 17.57     | 13.730     | 16.419     |
| 06-08-2006                     | 08:41:25 | 13.325  | 18.12     | 9.852      | 16.230     |
| 06-08-2006                     | 08:42:25 | 13.460  | 19.04     | 8.391      | 16.108     |
| 06-08-2006                     | 08:43:24 | 13.473  | 20.13     | 11.443     | 15.797     |
| 06-08-2006                     | 08:44:25 | 13.531  | 19.32     | 6.173      | 15.743     |
| 06-08-2006                     | 08:45:25 | 13.405  | 21.13     | 5.781      | 16.021     |
| 06-08-2006                     | 08:46:24 | 13.453  | 19.05     | 4.610      | 15.912     |
| 06-08-2006                     | 08:47:25 | 13.232  | 16.46     | 3.491      | 16.580     |
| 06-08-2006                     | 08:48:24 | 13.149  | 17.27     | 3.507      | 16.560     |
| 06-08-2006                     | 08:49:25 | 13.303  | 20.21     | 4.434      | 16.162     |
| 06-08-2006                     | 08:50:25 | 13.282  | 18.95     | 4.121      | 16.163     |
| 06-08-2006                     | 08:51:24 | 13.216  | 19.05     | 3.618      | 16.338     |
| Begin calculating run averages |          |         |           |            |            |
| 06-08-2006                     | 08:52:25 | 13.099  | 16.25     | 2.739      | 16.618     |
| 06-08-2006                     | 08:53:26 | 13.032  | 14.85     | 2.337      | 16.777     |
| 06-08-2006                     | 08:54:25 | 12.961  | 13.80     | 2.387      | 16.793     |
| 06-08-2006                     | 08:55:25 | 13.157  | 16.04     | 2.755      | 16.575     |
| 06-08-2006                     | 08:56:24 | 13.136  | 15.79     | 2.905      | 16.427     |
| 06-08-2006                     | 08:57:25 | 13.132  | 14.05     | 2.782      | 16.729     |
| 06-08-2006                     | 08:58:25 | 12.768  | 10.70     | 1.775      | 17.398     |
| 06-08-2006                     | 08:59:24 | 13.071  | 15.53     | 2.935      | 16.494     |
| 06-08-2006                     | 09:00:25 | 13.140  | 16.10     | 3.277      | 16.649     |
| 06-08-2006                     | 09:01:26 | 13.027  | 15.59     | 2.775      | 16.846     |
| 06-08-2006                     | 09:02:25 | 13.249  | 16.42     | 3.163      | 16.276     |
| 06-08-2006                     | 09:03:25 | 13.135  | 16.11     | 2.972      | 16.689     |
| 06-08-2006                     | 09:04:24 | 13.182  | 15.81     | 2.784      | 16.456     |
| 06-08-2006                     | 09:05:25 | 13.285  | 17.45     | 3.077      | 16.295     |
| 06-08-2006                     | 09:06:25 | 13.239  | 15.87     | 2.647      | 16.308     |
| 06-08-2006                     | 09:07:24 | 13.216  | 14.90     | 2.529      | 16.717     |
| 06-08-2006                     | 09:08:25 | 13.123  | 12.59     | 2.259      | 16.541     |
| 06-08-2006                     | 09:09:26 | 13.116  | 13.47     | 2.089      | 16.900     |
| 06-08-2006                     | 09:10:25 | 12.535  | 8.34      | 0.928      | 18.339     |
| 06-08-2006                     | 09:11:25 | 12.521  | 9.04      | 0.829      | 18.208     |
| 06-08-2006                     | 09:12:24 | 12.322  | 7.06      | 0.615      | 18.752     |
| 06-08-2006                     | 09:13:25 | 12.884  | 14.30     | 1.796      | 17.029     |
| 06-08-2006                     | 09:14:25 | 12.760  | 10.90     | 1.217      | 17.662     |
| 06-08-2006                     | 09:15:24 | 12.846  | 11.50     | 1.357      | 17.421     |
| 06-08-2006                     | 09:16:25 | 12.830  | 12.02     | 1.412      | 17.387     |
| 06-08-2006                     | 09:17:26 | 13.075  | 16.56     | 2.444      | 16.523     |

Test Run 1 STRATA Version 1.1

|                     | O2<br>% | CO<br>ppm | THC<br>ppm | NOx<br>ppm |
|---------------------|---------|-----------|------------|------------|
| 06-08-2006 09:18:25 | 13.059  | 14.21     | 1.885      | 16.812     |
| 06-08-2006 09:19:25 | 13.204  | 14.91     | 2.136      | 16.523     |
| 06-08-2006 09:20:24 | 12.789  | 9.31      | 1.056      | 17.841     |
| 06-08-2006 09:21:25 | 12.863  | 11.56     | 1.310      | 17.116     |
| Run Averages        | O2<br>% | CO<br>ppm | THC<br>ppm | NOx<br>ppm |
| 06-08-2006 09:22:05 | 12.994  | 13.75     | 2.175      | 16.970     |

Operator: B Johnston  
 Plant Name: Redwood Landfill  
 Location: Flare A51  
 Test Run 1 End

Final System Bias Check, Run 1 STRATA Version 1.1

|                     | O2<br>% | CO<br>ppm | THC<br>ppm | NOx<br>ppm |
|---------------------|---------|-----------|------------|------------|
| 06-08-2006 09:23:22 | 4.403   | 2.64      | 0.286      | 1.165      |
| 06-08-2006 09:24:23 | 0.167   | 0.86      | -0.082     | 0.006      |
| 06-08-2006 09:25:21 | 0.092   | -0.15     | -0.082     | 0.006      |
| 06-08-2006 09:26:22 | 0.075   | -0.16     | -0.083     | 0.006      |
| 06-08-2006 09:27:21 | 13.577  | 75.37     | 45.423     | 37.758     |
| 06-08-2006 09:28:22 | 18.865  | 91.01     | 47.032     | 44.324     |
| 06-08-2006 09:29:22 | 18.954  | 91.07     | 44.992     | 44.682     |
| 06-08-2006 09:30:21 | 18.966  | 91.16     | 44.920     | 44.700     |

Final System Bias Check for Run 1

Operator: B Johnston  
 Plant Name: Redwood Landfill  
 Location: Flare A51

Reference Cylinder Numbers

|     | Zero | Span     |
|-----|------|----------|
| O2  |      | CC8445   |
| CO  |      | AAL19769 |
| THC |      | AAL4576  |
| NOx |      | AAL19769 |

| Date/Time             | 06-08-2006 | 09:30:35 | PASSED |
|-----------------------|------------|----------|--------|
| Analyte               | O2         | CO       | THC    |
| Units                 | %          | ppm      | ppm    |
| Zero Ref Cyl          | 0.000      | 0.00     | 0.000  |
| Zero Cal              | 0.000      | 0.00     | 0.000  |
| Zero Avg              | 0.076      | -0.13    | -0.083 |
| Zero Bias%            | 0.3%       | 0.1%     | 0.2%   |
| Zero Drift%           | -0.1%      | 0.0%     | 0.0%   |
| Span Ref Cyl          | 19.040     | 90.10    | 45.000 |
| Span Cal              | 19.040     | 90.10    | 45.000 |
| Span Avg              | 18.967     | 91.21    | 44.908 |
| Span Bias%            | 0.3%       | 1.1%     | 0.2%   |
| Span Drift%           | -0.5%      | 0.7%     | -0.4%  |
| Ini Zero Avg          | 0.105      | -0.10    | -0.076 |
| Ini Span Avg          | 19.088     | 90.46    | 45.090 |
| Run Avg               | 12.994     | 13.75    | 2.175  |
| Co                    | 0.091      | -0.11    | -0.079 |
| Cm                    | 19.028     | 90.83    | 44.999 |
| Correct Avg           | 12.973     | 13.73    | 2.251  |
| System Bias Check End |            |          |        |

Test Run 2 STRATA Version 1.1

|                                | O2<br>% | CO<br>ppm | THC<br>ppm | NOx<br>ppm |
|--------------------------------|---------|-----------|------------|------------|
| 06-08-2006 09:31:36            | 15.152  | 17.75     | 9.498      | 20.581     |
| 06-08-2006 09:32:36            | 12.771  | 9.07      | 0.742      | 17.750     |
| Begin calculating run averages |         |           |            |            |
| 06-08-2006 09:33:35            | 12.529  | 6.54      | 0.413      | 18.329     |
| 06-08-2006 09:34:36            | 12.478  | 7.09      | 0.480      | 18.301     |
| 06-08-2006 09:35:35            | 12.204  | 5.00      | 0.224      | 18.880     |
| 06-08-2006 09:36:35            | 12.258  | 5.33      | 0.266      | 18.687     |
| 06-08-2006 09:37:36            | 12.437  | 10.96     | 0.832      | 18.013     |
| 06-08-2006 09:38:35            | 12.844  | 13.22     | 1.326      | 16.756     |
| 06-08-2006 09:39:36            | 12.876  | 10.52     | 1.244      | 17.155     |
| 06-08-2006 09:40:36            | 13.010  | 16.24     | 2.206      | 16.396     |
| 06-08-2006 09:41:35            | 12.832  | 9.45      | 1.014      | 17.622     |
| 06-08-2006 09:42:36            | 13.000  | 13.23     | 1.580      | 16.460     |
| 06-08-2006 09:43:35            | 13.022  | 13.34     | 1.509      | 16.565     |
| 06-08-2006 09:44:35            | 13.089  | 12.98     | 1.785      | 16.248     |
| 06-08-2006 09:45:36            | 13.179  | 13.83     | 1.856      | 16.277     |
| 06-08-2006 09:46:35            | 13.046  | 13.19     | 1.515      | 16.594     |
| 06-08-2006 09:47:36            | 13.254  | 14.89     | 2.141      | 15.857     |
| 06-08-2006 09:48:36            | 13.130  | 14.82     | 1.912      | 16.351     |
| 06-08-2006 09:49:35            | 12.843  | 10.99     | 1.127      | 17.201     |
| 06-08-2006 09:50:36            | 12.598  | 10.44     | 1.085      | 17.537     |
| 06-08-2006 09:51:35            | 12.867  | 15.42     | 1.855      | 16.774     |
| 06-08-2006 09:52:35            | 12.739  | 15.11     | 1.840      | 17.222     |
| 06-08-2006 09:53:36            | 13.055  | 14.08     | 1.520      | 16.346     |
| 06-08-2006 09:54:35            | 13.032  | 13.38     | 1.525      | 16.434     |
| 06-08-2006 09:55:36            | 12.696  | 9.22      | 0.736      | 17.521     |
| 06-08-2006 09:56:36            | 12.961  | 12.93     | 1.450      | 16.561     |
| 06-08-2006 09:57:35            | 12.902  | 10.57     | 1.058      | 16.859     |
| 06-08-2006 09:58:36            | 12.823  | 10.47     | 1.139      | 17.112     |
| 06-08-2006 09:59:35            | 12.818  | 10.85     | 1.179      | 17.131     |
| 06-08-2006 10:00:35            | 12.594  | 9.17      | 0.769      | 17.716     |
| 06-08-2006 10:01:36            | 12.516  | 10.41     | 0.755      | 17.490     |
| 06-08-2006 10:02:35            | 12.777  | 13.71     | 1.178      | 16.967     |
| Run Averages                   | O2<br>% | CO<br>ppm | THC<br>ppm | NOx<br>ppm |
| 06-08-2006 10:03:04            | 12.815  | 11.65     | 1.257      | 17.095     |

Operator: B Johnston  
Plant Name: Redwood Landfill  
Location: Flare A51  
Test Run 2 End

# Final System Bias Check, Run 2 STRATA Version 1.1

|                     | O2<br>% | CO<br>ppm | THC<br>ppm | NOx<br>ppm |
|---------------------|---------|-----------|------------|------------|
| 06-08-2006 10:04:09 | 4.267   | 2.13      | 0.055      | 1.337      |
| 06-08-2006 10:05:10 | 0.141   | 1.00      | -0.088     | 0.002      |
| 06-08-2006 10:06:11 | 0.067   | 0.15      | -0.088     | 0.001      |
| 06-08-2006 10:07:10 | 8.895   | 53.98     | 29.886     | 28.494     |
| 06-08-2006 10:08:10 | 18.688  | 85.57     | 44.865     | 44.190     |
| 06-08-2006 10:09:09 | 18.870  | 85.70     | 44.820     | 44.363     |
| 06-08-2006 10:10:10 | 18.891  | 89.00     | 44.759     | 44.368     |
| 06-08-2006 10:11:11 | 18.897  | 89.65     | 44.725     | 44.372     |

## Final System Bias Check for Run 2

Operator: B Johnston  
Plant Name: Redwood Landfill  
Location: Flare A51

### Reference Cylinder Numbers

|     | Zero | Span     |
|-----|------|----------|
| O2  |      | CC8445   |
| CO  |      | AAL19769 |
| THC |      | AAL4576  |
| NOx |      | AAL19769 |

| Date/Time             | 06-08-2006 | 10:11:35 | PASSED |
|-----------------------|------------|----------|--------|
| Analyte               | O2         | CO       | THC    |
| Units                 | %          | ppm      | ppm    |
| Zero Ref Cyl          | 0.000      | 0.00     | 0.000  |
| Zero Cal              | 0.000      | 0.00     | 0.000  |
| Zero Avg              | 0.063      | 0.15     | -0.088 |
| Zero Bias%            | 0.3%       | 0.1%     | 0.2%   |
| Zero Drift%           | -0.1%      | 0.3%     | 0.0%   |
| Span Ref Cyl          | 19.040     | 90.10    | 45.000 |
| Span Cal              | 19.040     | 90.10    | 45.000 |
| Span Avg              | 18.899     | 89.74    | 44.750 |
| Span Bias%            | 0.6%       | 0.4%     | 0.5%   |
| Span Drift%           | -0.3%      | -1.5%    | -0.3%  |
| Ini Zero Avg          | 0.076      | -0.13    | -0.083 |
| Ini Span Avg          | 18.967     | 91.21    | 44.908 |
| Run Avg               | 12.815     | 11.65    | 1.257  |
| Co                    | 0.070      | 0.01     | -0.085 |
| Cm                    | 18.933     | 90.47    | 44.829 |
| Correct Avg           | 12.864     | 11.59    | 1.345  |
| System Bias Check End |            |          |        |

Test Run 3 STRATA Version 1.1

|                                | O2<br>% | CO<br>ppm | THC<br>ppm | NOx<br>ppm |
|--------------------------------|---------|-----------|------------|------------|
| 06-08-2006 10:12:36            | 15.099  | 36.53     | 7.025      | 20.686     |
| 06-08-2006 10:13:35            | 13.301  | 24.97     | 6.535      | 15.483     |
| Begin calculating run averages |         |           |            |            |
| 06-08-2006 10:14:36            | 13.640  | 28.20     | 8.525      | 15.153     |
| 06-08-2006 10:15:36            | 13.082  | 22.65     | 7.038      | 16.310     |
| 06-08-2006 10:16:35            | 13.312  | 24.11     | 6.875      | 15.383     |
| 06-08-2006 10:17:36            | 13.552  | 27.37     | 9.210      | 14.762     |
| 06-08-2006 10:18:36            | 13.594  | 28.37     | 7.801      | 14.754     |
| 06-08-2006 10:19:35            | 13.265  | 25.96     | 8.982      | 15.584     |
| 06-08-2006 10:20:36            | 13.365  | 21.96     | 5.193      | 15.145     |
| 06-08-2006 10:21:35            | 13.511  | 26.67     | 6.712      | 14.871     |
| 06-08-2006 10:22:36            | 13.463  | 26.50     | 6.799      | 15.140     |
| 06-08-2006 10:23:36            | 13.359  | 21.05     | 3.762      | 15.420     |
| 06-08-2006 10:24:35            | 13.165  | 18.41     | 2.853      | 15.573     |
| 06-08-2006 10:25:36            | 13.234  | 18.51     | 2.940      | 15.588     |
| 06-08-2006 10:26:37            | 13.067  | 16.75     | 2.261      | 16.088     |
| 06-08-2006 10:27:35            | 12.968  | 16.61     | 2.007      | 16.079     |
| 06-08-2006 10:28:36            | 13.002  | 15.89     | 1.612      | 16.130     |
| 06-08-2006 10:29:35            | 12.862  | 14.20     | 1.309      | 16.422     |
| 06-08-2006 10:30:36            | 12.898  | 15.75     | 1.471      | 16.387     |
| 06-08-2006 10:31:36            | 12.860  | 15.68     | 1.370      | 16.366     |
| 06-08-2006 10:32:35            | 12.804  | 15.05     | 1.186      | 16.474     |
| 06-08-2006 10:33:36            | 13.257  | 19.87     | 2.510      | 15.609     |
| 06-08-2006 10:34:37            | 13.094  | 21.49     | 3.316      | 16.435     |
| 06-08-2006 10:35:36            | 13.088  | 22.23     | 4.092      | 15.821     |
| 06-08-2006 10:36:36            | 13.529  | 25.25     | 4.815      | 14.748     |
| 06-08-2006 10:37:35            | 13.402  | 24.06     | 5.744      | 15.147     |
| 06-08-2006 10:38:36            | 13.284  | 21.20     | 4.999      | 16.035     |
| 06-08-2006 10:39:36            | 12.980  | 21.59     | 5.014      | 16.212     |
| 06-08-2006 10:40:35            | 12.999  | 21.14     | 6.563      | 16.142     |
| 06-08-2006 10:41:36            | 13.050  | 19.77     | 5.906      | 16.211     |
| 06-08-2006 10:42:35            | 13.215  | 22.34     | 8.585      | 15.361     |
| 06-08-2006 10:43:36            | 13.475  | 25.91     | 7.689      | 14.889     |
| Run Averages                   | O2<br>% | CO<br>ppm | THC<br>ppm | NOx<br>ppm |
| 06-08-2006 10:44:03            | 13.210  | 21.46     | 4.900      | 15.671     |

Operator: B Johnston  
Plant Name: Redwood Landfill  
Location: Flare A51  
Test Run 3 End

# Final System Bias Check, Run 3 STRATA Version 1.1

|                     | O2<br>% | CO<br>ppm | THC<br>ppm | NOx<br>ppm |
|---------------------|---------|-----------|------------|------------|
| 06-08-2006 10:45:25 | 3.707   | 3.31      | 0.925      | 0.396      |
| 06-08-2006 10:46:24 | 0.113   | 0.19      | -0.090     | -0.001     |
| 06-08-2006 10:47:24 | 0.051   | 0.00      | -0.091     | -0.001     |
| 06-08-2006 10:48:25 | 10.299  | 62.49     | 32.805     | 31.882     |
| 06-08-2006 10:49:24 | 18.737  | 90.68     | 44.428     | 35.776     |
| 06-08-2006 10:50:25 | 18.866  | 90.88     | 44.736     | 44.866     |
| 06-08-2006 10:51:24 | 18.883  | 90.57     | 44.717     | 44.869     |
| 06-08-2006 10:52:24 | 18.891  | 91.00     | 44.724     | 44.874     |

## Final System Bias Check for Run 3

Operator: B Johnston  
 Plant Name: Redwood Landfill  
 Location: Flare A51

### Reference Cylinder Numbers

|     | Zero | Span     |
|-----|------|----------|
| O2  |      | CC8445   |
| CO  |      | AAL19769 |
| THC |      | AAL4576  |
| NOx |      | AAL19769 |

| Date/Time             | 06-08-2006 | 10:52:37 | PASSED |
|-----------------------|------------|----------|--------|
| Analyte               | O2         | CO       | THC    |
| Units                 | %          | ppm      | ppm    |
| Zero Ref Cyl          | 0.000      | 0.00     | 0.000  |
| Zero Cal              | 0.000      | 0.00     | 0.000  |
| Zero Avg              | 0.051      | 0.02     | -0.091 |
| Zero Bias%            | 0.2%       | 0.0%     | 0.2%   |
| Zero Drift%           | 0.0%       | -0.1%    | 0.0%   |
| Span Ref Cyl          | 19.040     | 90.10    | 45.000 |
| Span Cal              | 19.040     | 90.10    | 45.000 |
| Span Avg              | 18.891     | 91.03    | 44.734 |
| Span Bias%            | 0.6%       | 0.9%     | 0.5%   |
| Span Drift%           | 0.0%       | 1.3%     | 0.0%   |
| Ini Zero Avg          | 0.063      | 0.15     | -0.088 |
| Ini Span Avg          | 18.899     | 89.74    | 44.750 |
| Run Avg               | 13.210     | 21.46    | 4.900  |
| Co                    | 0.057      | 0.08     | -0.089 |
| Cm                    | 18.895     | 90.39    | 44.742 |
| Correct Avg           | 13.294     | 21.33    | 5.007  |
| System Bias Check End |            |          |        |

PART NO. LG/8011

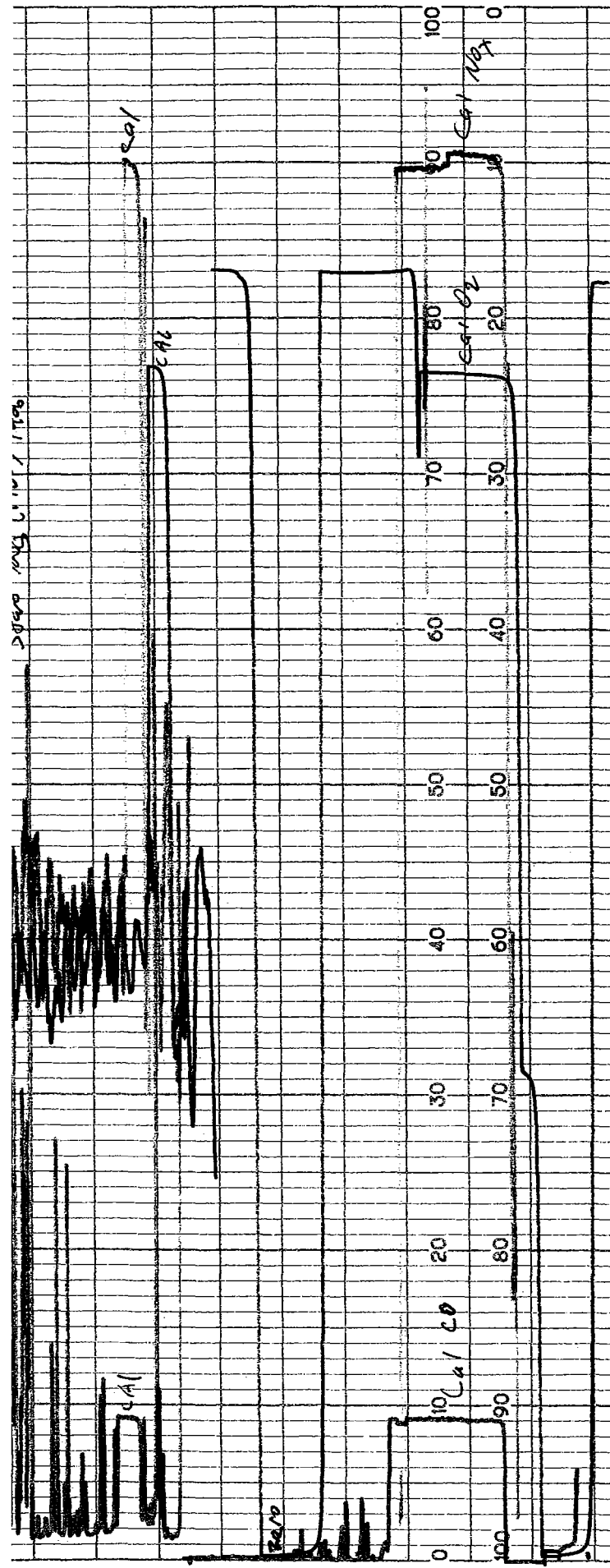
PRINTED IN U.S.A.

Blue: O<sub>2</sub> 0-25%  
Green: CO 0-1000 PPM  
Red: NO<sub>x</sub> 0-50 PPM

Pretest A-50

LINE 5 E 1 S

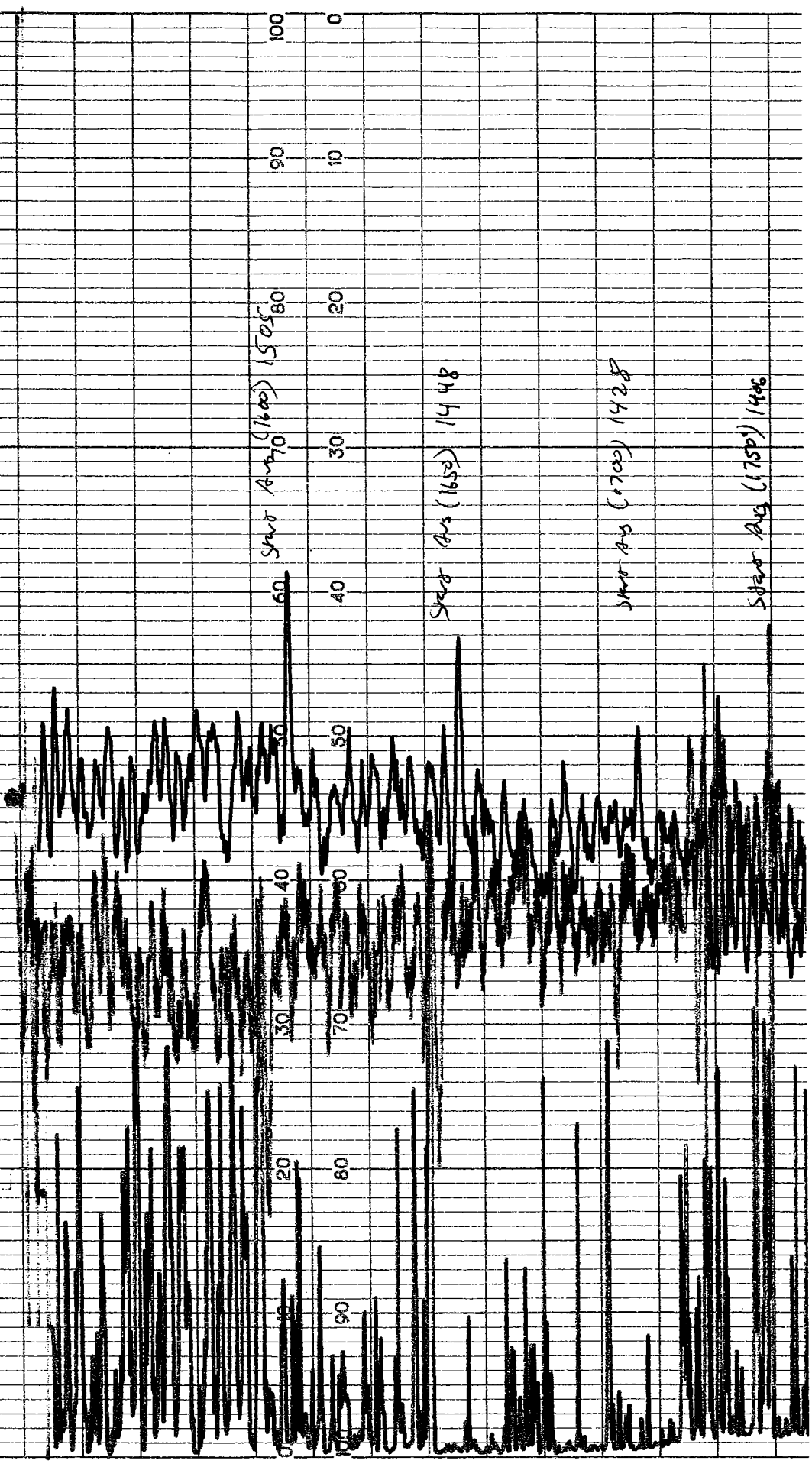
Redwood Landfill 6-6-06

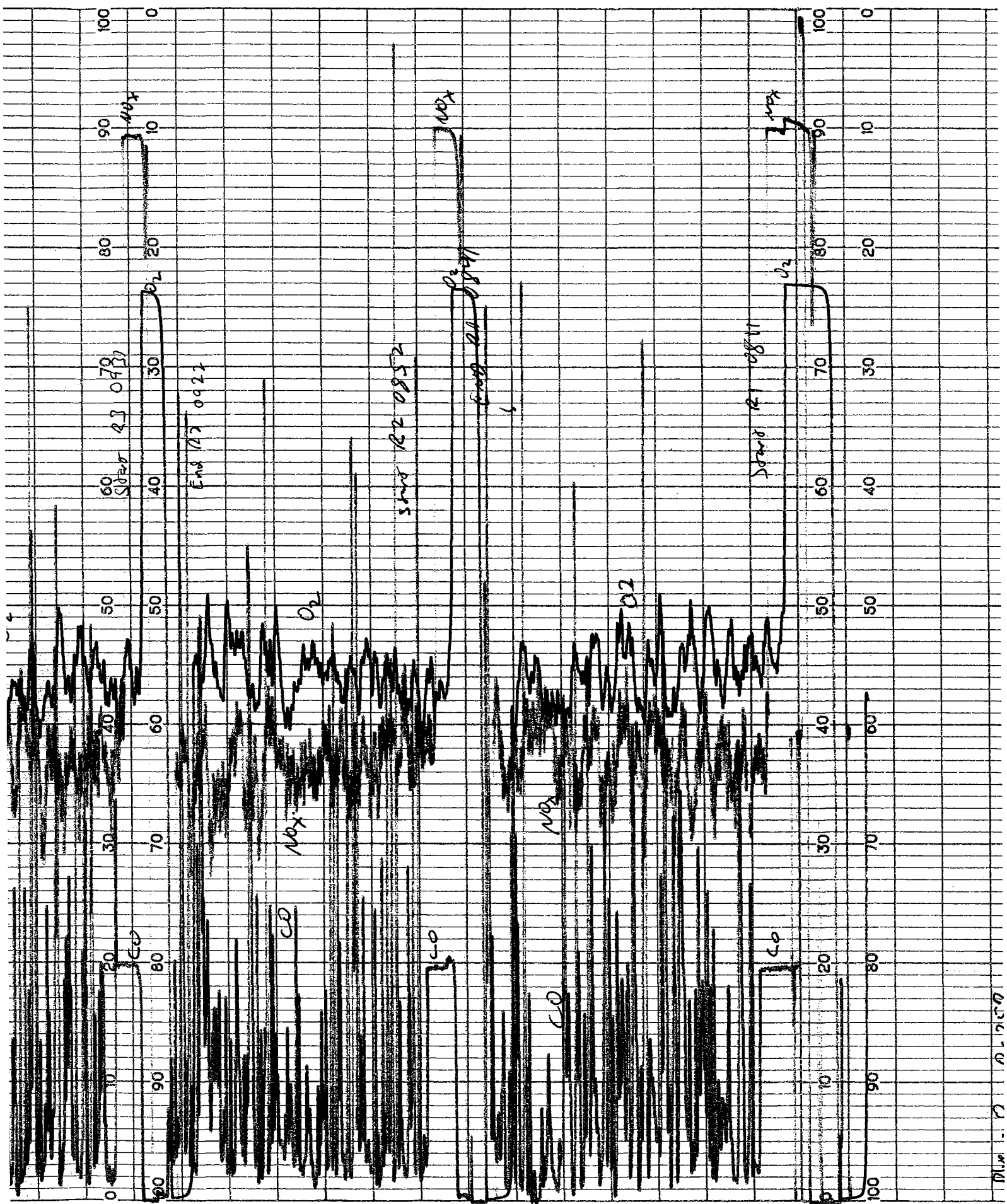


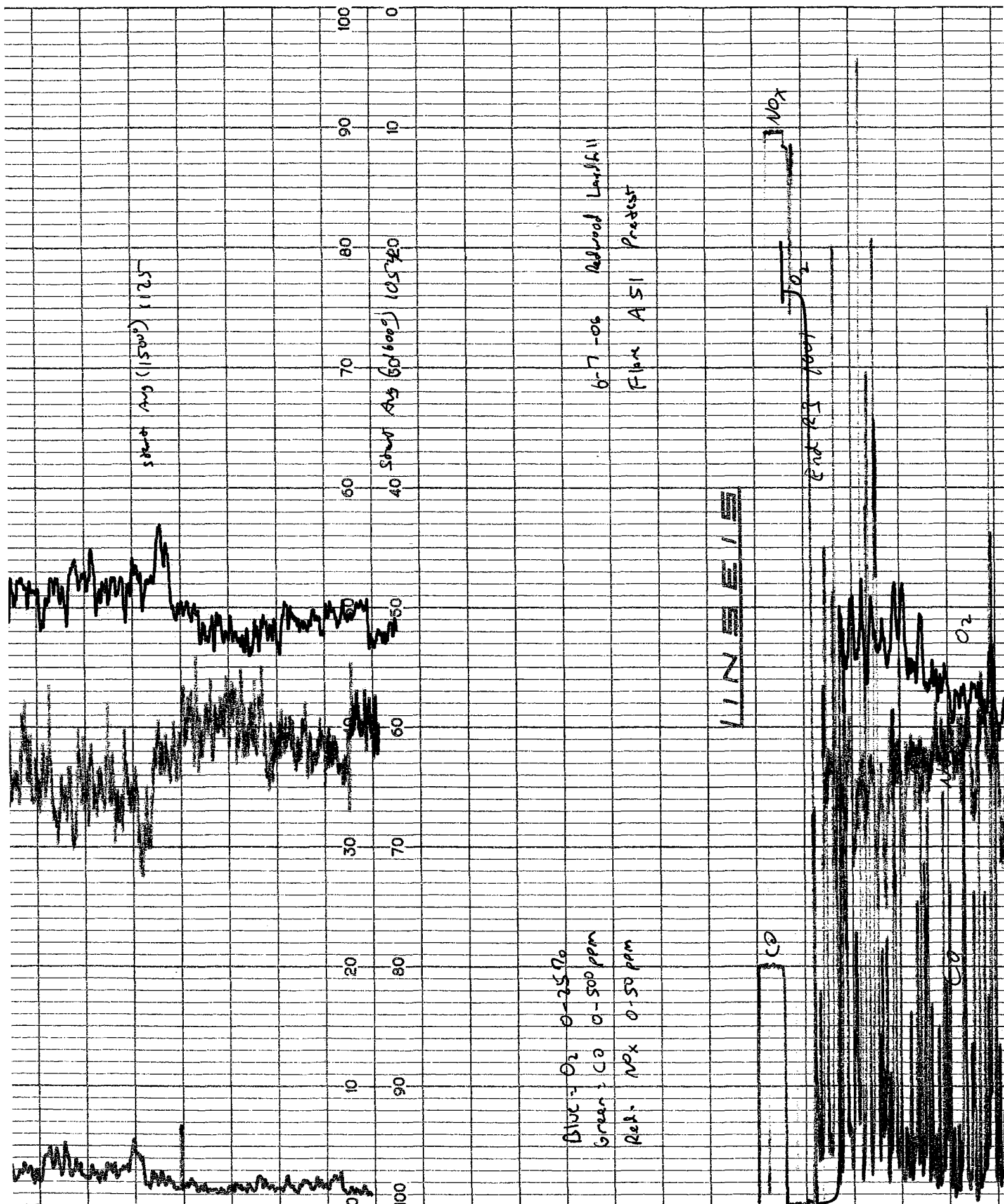
Blue =  $O_2$  0-25%  
Green = CO 0-500 ppm  
Red = NOx 0-50 ppm

LINE 1

6-7-98 Redwood Landfill  
Flare A-50







PART NO. LO/8011

PRINTED IN U.S.A.

Blue =  $O_2$  0-25%

Red. NO<sub>x</sub> 0-50 ppm

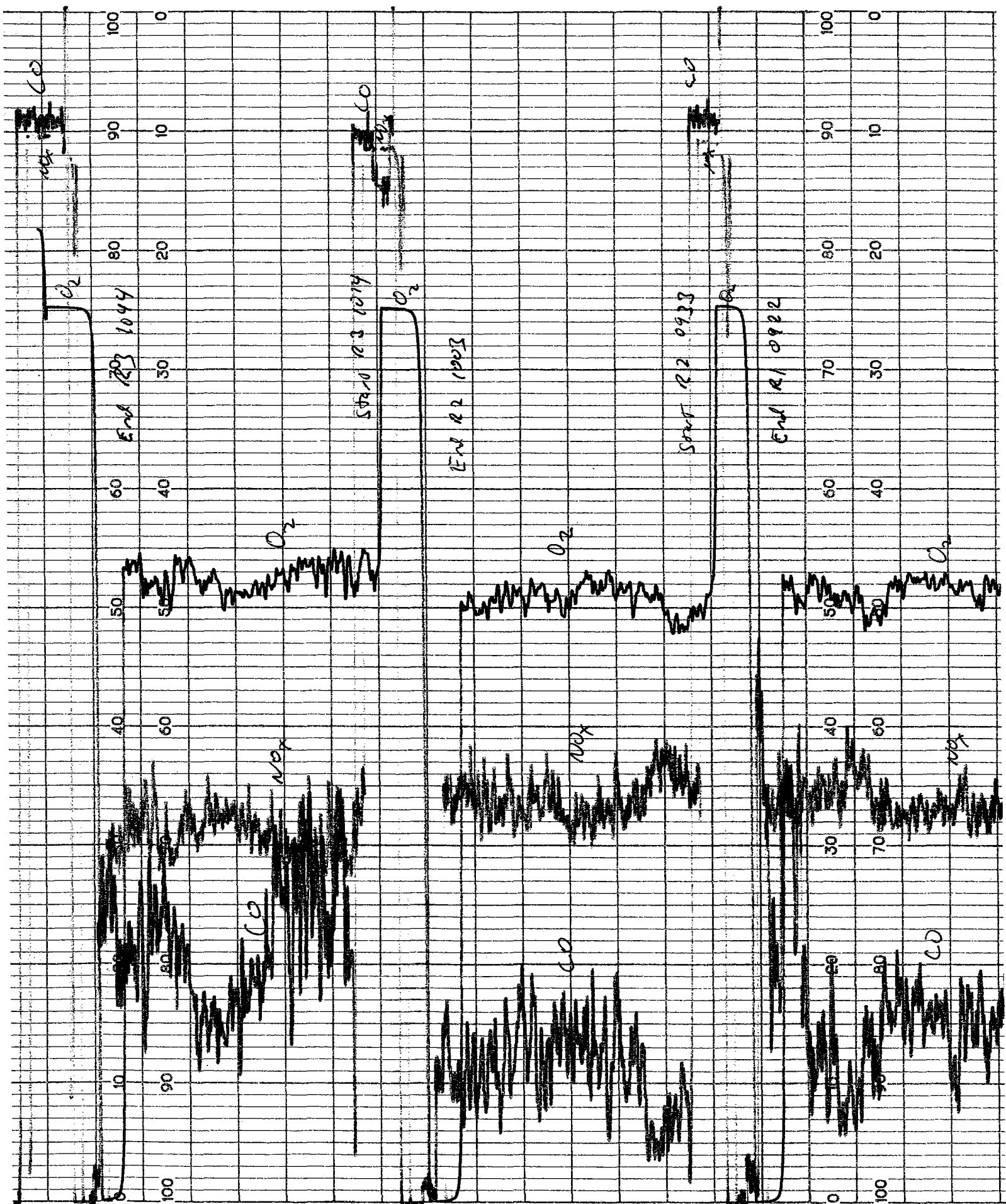
Q-20 PM

|            |         |           |
|------------|---------|-----------|
| 6-8-06     | Redwood | Land 6.11 |
| Elmore AS1 |         |           |

Flare A5-1

LS19 (0.0561) SNR 10.245

⑤



Star Avg (1450) 1448

Star Avg (1700) 1728

Star Avg (2800) 1406

(a)

Zero

Col FHE

Green : THC 0-50 ppm

Redwood Landfill 6-6-06  
Pretest A-50

Starts 21 08M

7.82

6-7-06 Redwood Landfill

Flare A-50

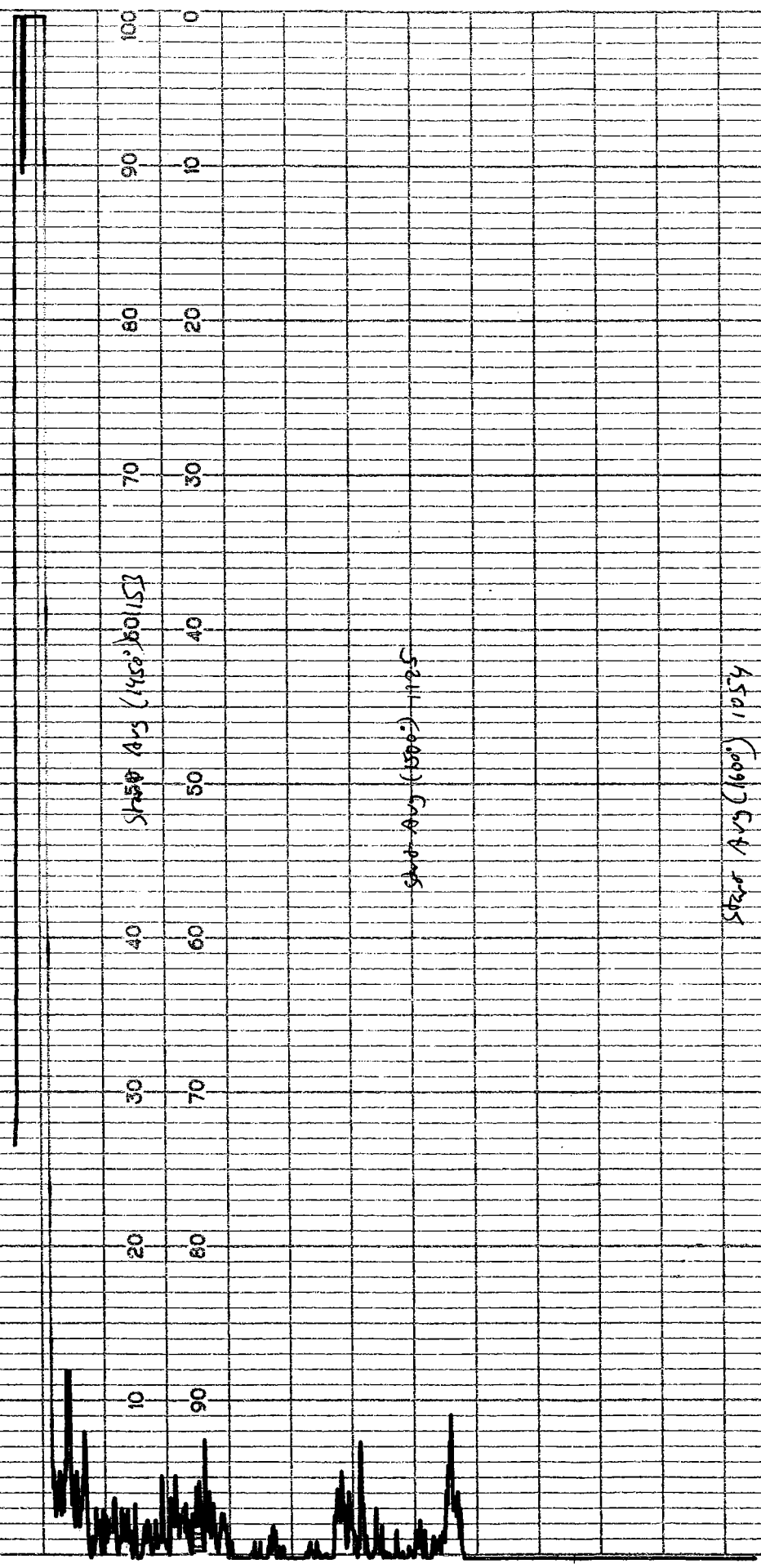
Green - TAC 0-50 PPM

Starts 21 08M

LINE 15



PRINTED



6-7-00 Railroad Carfill  
Flare A 51 Pre test

Green - T4C 0-50 ppm

5560 28 20955

End 0922

Steps R1 0852

Green = T11C 0-50 ppm

5-8-06 Redwood Landfill

Flare

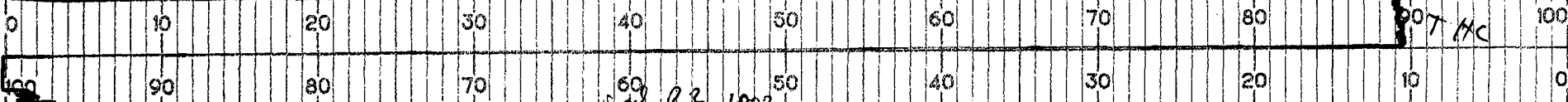
|                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                 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| 1. <b>RESEARCH</b><br>2. <b>DESIGN</b><br>3. <b>CONSTRUCTION</b><br>4. <b>OPERATION</b><br>5. <b>MAINTENANCE</b><br>6. <b>REPAIR</b><br>7. <b>REPLACEMENT</b><br>8. <b>RECONSTRUCTION</b><br>9. <b>RENOVATION</b><br>10. <b>REMODELING</b><br>11. <b>REPAIR</b><br>12. <b>REPLACEMENT</b><br>13. <b>RECONSTRUCTION</b><br>14. <b>RENOVATION</b><br>15. <b>REMODELING</b><br>16. <b>REPAIR</b><br>17. <b>REPLACEMENT</b><br>18. <b>RECONSTRUCTION</b><br>19. <b>RENOVATION</b><br>20. <b>REMODELING</b><br>21. <b>REPAIR</b><br>22. <b>REPLACEMENT</b><br>23. <b>RECONSTRUCTION</b><br>24. <b>RENOVATION</b><br>25. <b>REMODELING</b><br>26. <b>REPAIR</b><br>27. <b>REPLACEMENT</b><br>28. <b>RECONSTRUCTION</b><br>29. <b>RENOVATION</b><br>30. <b>REMODELING</b><br>31. <b>REPAIR</b><br>32. <b>REPLACEMENT</b><br>33. <b>RECONSTRUCTION</b><br>34. <b>RENOVATION</b><br>35. <b>REMODELING</b><br>36. <b>REPAIR</b><br>37. <b>REPLACEMENT</b><br>38. <b>RECONSTRUCTION</b><br>39. <b>RENOVATION</b><br>40. <b>REMODELING</b><br>41. <b>REPAIR</b><br>42. <b>REPLACEMENT</b><br>43. <b>RECONSTRUCTION</b><br>44. <b>RENOVATION</b><br>45. <b>REMODELING</b><br>46. <b>REPAIR</b><br>47. <b>REPLACEMENT</b><br>48. <b>RECONSTRUCTION</b><br>49. <b>RENOVATION</b><br>50. <b>REMODELING</b><br>51. <b>REPAIR</b><br>52. <b>REPLACEMENT</b><br>53. <b>RECONSTRUCTION</b><br>54. <b>RENOVATION</b><br>55. <b>REMODELING</b><br>56. <b>REPAIR</b><br>57. <b>REPLACEMENT</b><br>58. <b>RECONSTRUCTION</b><br>59. <b>RENOVATION</b><br>60. <b>REMODELING</b><br>61. <b>REPAIR</b><br>62. <b>REPLACEMENT</b><br>63. <b>RECONSTRUCTION</b><br>64. <b>RENOVATION</b><br>65. <b>REMODELING</b><br>66. <b>REPAIR</b><br>67. <b>REPLACEMENT</b><br>68. <b>RECONSTRUCTION</b><br>69. <b>RENOVATION</b><br>70. <b>REMODELING</b><br>71. <b>REPAIR</b><br>72. <b>REPLACEMENT</b><br>73. <b>RECONSTRUCTION</b><br>74. <b>RENOVATION</b><br>75. <b>REMODELING</b><br>76. <b>REPAIR</b><br>77. <b>REPLACEMENT</b><br>78. <b>RECONSTRUCTION</b><br>79. <b>RENOVATION</b><br>80. <b>REMODELING</b><br>81. <b>REPAIR</b><br>82. <b>REPLACEMENT</b><br>83. <b>RECONSTRUCTION</b><br>84. <b>RENOVATION</b><br>85. <b>REMODELING</b><br>86. <b>REPAIR</b><br>87. <b>REPLACEMENT</b><br>88. <b>RECONSTRUCTION</b><br>89. <b>RENOVATION</b><br>90. <b>REMODELING</b><br>91. <b>REPAIR</b><br>92. <b>REPLACEMENT</b><br>93. <b>RECONSTRUCTION</b><br>94. <b>RENOVATION</b><br>95. <b>REMODELING</b><br>96. <b>REPAIR</b><br>97. <b>REPLACEMENT</b><br>98. <b>RECONSTRUCTION</b><br>99. <b>RENOVATION</b><br>100. <b>REMODELING</b><br>101. <b>REPAIR</b><br>102. <b>REPLACEMENT</b><br>103. <b>RECONSTRUCTION</b><br>104. <b>RENOVATION</b><br>105. <b>REMODELING</b><br>106. <b>REPAIR</b><br>107. <b>REPLACEMENT</b><br>108. <b>RECONSTRUCTION</b><br>109. <b>RENOVATION</b><br>110. <b>REMODELING</b><br>111. <b>REPAIR</b><br>112. <b>REPLACEMENT</b><br>113. <b>RECONSTRUCTION</b><br>114. <b>RENOVATION</b><br>115. <b>REMODELING</b><br>116. <b>REPAIR</b><br>117. <b>REPLACEMENT</b><br>118. <b>RECONSTRUCTION</b><br>119. <b>RENOVATION</b><br>120. <b>REMODELING</b><br>121. <b>REPAIR</b><br>122. <b>REPLACEMENT</b><br>123. <b>RECONSTRUCTION</b><br>124. <b>RENOVATION</b><br>125. <b>REMODELING</b><br>126. <b>REPAIR</b><br>127. <b>REPLACEMENT</b><br>128. <b>RECONSTRUCTION</b><br>129. <b>RENOVATION</b><br>130. <b>REMODELING</b><br>131. <b>REPAIR</b><br>132. <b>REPLACEMENT</b><br>133. <b>RECONSTRUCTION</b><br>134. <b>RENOVATION</b><br>135. <b>REMODELING</b><br>136. <b>REPAIR</b><br>137. <b>REPLACEMENT</b><br>138. <b>RECONSTRUCTION</b><br>139. <b>RENOVATION</b><br>140. <b>REMODELING</b><br>141. <b>REPAIR</b><br>142. <b>REPLACEMENT</b><br>143. <b>RECONSTRUCTION</b><br>144. <b>RENOVATION</b><br>145. <b>REMODELING</b><br>146. <b>REPAIR</b><br>147. <b>REPLACEMENT</b><br>148. <b>RECONSTRUCTION</b><br>149. <b>RENOVATION</b><br>150. <b>REMODELING</b><br>151. <b>REPAIR</b><br>152. <b>REPLACEMENT</b><br>153. <b>RECONSTRUCTION</b><br>154. <b>RENOVATION</b><br>155. <b>REMODELING</b><br>156. <b>REPAIR</b><br>157. <b>REPLACEMENT</b><br>158. <b>RECONSTRUCTION</b><br>159. <b>RENOVATION</b><br>160. <b>REMODELING</b><br>161. <b>REPAIR</b><br>162. <b>REPLACEMENT</b><br>163. <b>RECONSTRUCTION</b><br>164. <b>RENOVATION</b><br>165. <b>REMODELING</b><br>166. <b>REPAIR</b><br>167. <b>REPLACEMENT</b><br>168. <b>RECONSTRUCTION</b><br>169. <b>RENOVATION</b><br>170. <b>REMODELING</b><br>171. <b>REPAIR</b><br>172. <b>REPLACEMENT</b><br>173. <b>RECONSTRUCTION</b><br>174. <b>RENOVATION</b><br>175. <b>REMODELING</b><br>176. <b>REPAIR</b><br>177. <b>REPLACEMENT</b><br>178. <b>RECONSTRUCTION</b><br>179. <b>RENOVATION</b><br>180. <b>REMODELING</b><br>181. <b>REPAIR</b><br>182. <b>REPLACEMENT</b><br>183. <b>RECONSTRUCTION</b><br>184. <b>RENOVATION</b><br>185. <b>REMODELING</b><br>186. <b>REPAIR</b><br>187. <b>REPLACEMENT</b><br>188. <b>RECONSTRUCTION</b><br>189. <b>RENOVATION</b><br>190. <b>REMODELING</b><br>191. <b>REPAIR</b><br>192. <b>REPLACEMENT</b><br>193. <b>RECONSTRUCTION</b><br>194. <b>RENOVATION</b><br>195. <b>REMODELING</b><br>196. <b>REPAIR</b><br>197. <b>REPLACEMENT</b><br>198. <b>RECONSTRUCTION</b><br>199. <b>RENOVATION</b><br>200. <b>REMODELING</b><br>201. <b>REPAIR</b><br>202. <b>REPLACEMENT</b><br>203. <b>RECONSTRUCTION</b><br>204. <b>RENOVATION</b><br>205. <b>REMODELING</b><br>206. <b>REPAIR</b><br>207. <b>REPLACEMENT</b><br>208. <b>RECONSTRUCTION</b><br>209. <b>RENOVATION</b><br>210. <b>REMODELING</b><br>211. <b>REPAIR</b><br>212. <b>REPLACEMENT</b><br>213. <b>RECONSTRUCTION</b><br>214. <b>RENOVATION</b><br>215. <b>REMODELING</b><br>216. <b>REPAIR</b><br>217. <b>REPLACEMENT</b><br>218. <b>RECONSTRUCTION</b><br>219. <b>RENOVATION</b><br>220. <b>REMODELING</b><br>221. <b>REPAIR</b><br>222. <b>REPLACEMENT</b><br>223. <b>RECONSTRUCTION</b><br>224. <b>RENOVATION</b><br>225. <b>REMODELING</b><br>226. <b>REPAIR</b><br>227. <b>REPLACEMENT</b><br>228. <b>RECONSTRUCTION</b><br>229. <b>RENOVATION</b><br>230. <b>REMODELING</b><br>231. <b>REPAIR</b><br>232. <b>REPLACEMENT</b><br>233. <b>RECONSTRUCTION</b><br> |
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LINE 1

End R3 1044

THC

Start R3 1014



THC

Start R2 0933

THC

End R1 0922

THC

Start R1 0852



# Scott Specialty Gases

500 WEAVER PARK RD, LONGMONT, CO 80501

## RATA CLASS

### Dual-Analyzed Calibration Standard

Phone: 888-253-1635

Fax: 303-772-7675

## CERTIFICATE OF ACCURACY: Interference Free™ Multi-Component EPA Protocol Gas

### Assay Laboratory

SCOTT SPECIALTY GASES  
500 WEAVER PARK RD  
LONGMONT, CO 80501

P.O. No.: 101520  
Project No.: 08-36059-001

### Customer

SHAW ENVIRONMENTAL & INFRASTRUCTURE

BILL JOHNSTON  
2360 BERING DRIVE  
SAN JOSE CA 95131

### ANALYTICAL INFORMATION

This certification was performed according to EPA Traceability Protocol For Assay & Certification of Gaseous Calibration Standards; Procedure G-1; September, 1997.

Cylinder Number: **AAL19769** Certification Date: **10May2006** Exp. Date: **09May2008**  
Cylinder Pressure\*\*\*: **1870 PSIG**

| COMPONENT                | CERTIFIED CONCENTRATION (Moles) | ACCURACY** | TRACEABILITY         |
|--------------------------|---------------------------------|------------|----------------------|
| CARBON MONOXIDE          | 90.1 PPM                        | +/- 1%     | Direct NIST and NMI  |
| NITRIC OXIDE             | 44.7 PPM                        | +/- 1%     | Direct NIST and NMI  |
| NITROGEN - OXYGEN FREE   | BALANCE                         |            |                      |
| TOTAL OXIDES OF NITROGEN | 44.8 PPM                        |            | Reference Value Only |

\*\*\* Do not use when cylinder pressure is below 150 psig.

\*\* Analytical accuracy is based on the requirements of EPA Protocol Procedure G1, September 1997.

### REFERENCE STANDARD

| TYPE/SRM NO. | EXPIRATION DATE | CYLINDER NUMBER | CONCENTRATION | COMPONENT       |
|--------------|-----------------|-----------------|---------------|-----------------|
| NTRM 1679    | 02Apr2007       | ALM030197       | 94.90 PPM     | CARBON MONOXIDE |
| NTRM 1683    | 15Aug2009       | AAL070639       | 49.82 PPM     | NITRIC OXIDE    |

### INSTRUMENTATION

| INSTRUMENT/MODEL/SERIAL# | DATE LAST CALIBRATED | ANALYTICAL PRINCIPLE |
|--------------------------|----------------------|----------------------|
| FTIR//000929062          | 10May2006            | FTIR                 |
| FTIR//000929062          | 08May2006            | FTIR                 |

### ANALYZER READINGS

(Z=Zero Gas R=Reference Gas T=Test Gas r=Correlation Coefficient)

#### First Triad Analysis

#### CARBON MONOXIDE

Date: 01May2006 Response Unit: PPM  
Z1=-0.01877 R1=95.10122 T1=90.59397  
R2=95.12704 Z2=0.00010 T2=90.60740  
Z3=0.02836 T3=90.63578 R3=95.16162  
Avg. Concentration: 90.39 PPM

#### Second Triad Analysis

Date: 10May2006 Response Unit: PPM  
Z1=-0.02533 R1=95.68689 T1=90.52361  
R2=95.79857 Z2=-0.01272 T2=90.56149  
Z3=0.03944 T3=90.72761 R3=95.90278  
Avg. Concentration: 89.76 PPM

#### Calibration Curve

Concentration = A + Bx + Cx<sup>2</sup> + Dx<sup>3</sup> + Ex<sup>4</sup>  
r = 9.99985E-1  
Constants: A = 0.00000E+0  
B = 7.18007E-1 C = 3.18000E-4  
D = 0.00000E+0 E = 0.00000E+0

#### NITRIC OXIDE

Date: 01May2006 Response Unit: PPM  
Z1=-0.02773 R1=49.51237 T1=44.39455  
R2=49.69193 Z2=-0.01858 T2=44.81024  
Z3=0.11894 T3=44.90252 R3=49.72640  
Avg. Concentration: 44.86 PPM

Date: 10May2006 Response Unit: PPM  
Z1=-0.06655 R1=50.11168 T1=44.73440  
R2=50.12921 Z2=-0.05510 T2=44.78113  
Z3=-0.03373 T3=44.85699 R3=50.22003  
Avg. Concentration: 44.50 PPM

Concentration = A + Bx + Cx<sup>2</sup> + Dx<sup>3</sup> + Ex<sup>4</sup>  
r = 9.99993E-1  
Constants: A = 0.00000E+0  
B = 9.79296E-1 C = 1.73000E-4  
D = 0.00000E+0 E = 0.00000E+0

Special Notes:

PLEASE REPORT NOX

APPROVED BY:

Jon Witzak



**Scott Specialty Gases**

500 WEAVER PARK RD, LONGMONT, CO 80501

**CERTIFIED MASTER CLASS**

*Single-Certified Calibration Standard*

Phone: 888-253-1635

Fax: 303-772-7673

**CERTIFICATE OF ACCURACY: Certified Master Class Calibration Standard**

**Product Information**

Project No.: 08-24834-001

Item No.: 08022711 PAL

P.O. No.: SUSAN M. POWERS

Cylinder Number: AAL4576

Cylinder Size: AL

Certification Date: 25May2005

Expiration Date: 25May2008

**Customer**

SHAW ENVIRONMENTAL & INFRASTRUCTURE

2360 BERING DRIVE

SAN JOSE, CA 95131

**CERTIFIED CONCENTRATION**

**Component Name**

**Concentration  
(Moles)**

**Accuracy  
(+/-%)**

METHANE  
AIR

45.0 PPM  
BALANCE

2

**TRACEABILITY**

**Traceable To**

NIST

APPROVED BY:

  
JON WITZAK

DATE:

5-25-05



Praxair  
5700 South Alameda Street  
Los Angeles, CA 90058  
Telephone: (323) 585-2154  
Facsimile: (714) 542-6689

## CERTIFICATE OF ANALYSIS / EPA PROTOCOL GAS

CUSTOMER SHAW ENV.

P.O NUMBER

### REFERENCE STANDARD

| COMPONENT           | NIST SRM NO. | CYLINDER NO. | CONCENTRATION |
|---------------------|--------------|--------------|---------------|
| CARBON DIOXIDE GMIS | vs. SRM#2745 | CC 102004    | 19.73 %       |
| OXYGEN GMIS         | vs. SRM#2659 | CC 95713     | 21.04 %       |

### ANALYZER READINGS

R=REFERENCE STANDARD

Z=ZERO GAS

C=GAS CANDIDATE

|                      |                |                 |                         |                     |             |                 |   |       |
|----------------------|----------------|-----------------|-------------------------|---------------------|-------------|-----------------|---|-------|
| 1. COMPONENT         | CARBON DIOXIDE | GMIS            | ANALYZER MAKE-MODEL-S/N | Siemens Ultramat 5E | S/N A12-730 |                 |   |       |
| ANALYTICAL PRINCIPLE | NDIR           |                 | LAST CALIBRATION DATE   | 08/01/05            |             |                 |   |       |
| FIRST ANALYSIS DATE  | 08/08/05       |                 | SECOND ANALYSIS DATE    |                     |             |                 |   |       |
| Z 0.00               | R 19.74        | C 17.82         | CONC.                   | 17.81               | Z           | R               | C | CONC. |
| R 19.74              | Z 0.00         | C 17.80         | CONC.                   | 17.81               | R           | Z               | C | CONC. |
| Z 0.00               | C 17.80        | R 19.74         | CONC.                   | 17.81               | Z           | C               | R | CONC. |
| U/M %                |                | MEAN TEST ASSAY | 17.81                   | U/M %               |             | MEAN TEST ASSAY |   |       |
| 2. COMPONENT         | OXYGEN         | GMIS            | ANALYZER MAKE-MODEL-S/N | Siemens Oxymat 5E   | S/N A12-839 |                 |   |       |
| ANALYTICAL PRINCIPLE | Paramagnetic   |                 | LAST CALIBRATION DATE   | 08/01/05            |             |                 |   |       |
| FIRST ANALYSIS DATE  | 08/08/05       |                 | SECOND ANALYSIS DATE    |                     |             |                 |   |       |
| Z 0.00               | R 21.04        | C 19.04         | CONC.                   | 19.04               | Z           | R               | C | CONC. |
| R 21.04              | Z 0.00         | C 19.04         | CONC.                   | 19.04               | R           | Z               | C | CONC. |
| Z 0.00               | C 19.04        | R 21.04         | CONC.                   | 19.04               | Z           | C               | R | CONC. |
| U/M %                |                | MEAN TEST ASSAY | 19.04                   | U/M %               |             | MEAN TEST ASSAY |   |       |

VALUES NOT VALID BELOW 150 PSIG.

|                                         |                      |                         |           |
|-----------------------------------------|----------------------|-------------------------|-----------|
| THIS CYLINDER NO.                       | CC 8445              | CERTIFIED CONCENTRATION |           |
| HAS BEEN CERTIFIED ACCORDING TO SECTION | EPA-600/R97/121      | CARBON DIOXIDE          | 17.81 %   |
| OF TRACEABILITY PROTOCOL NO.            | REV 9/97             | OXYGEN                  | 19.04 %   |
| PROCEDURE                               | G1                   | NITROGEN                | BALANCE   |
| CERTIFIED ACCURACY                      | ± 1 % NIST TRACEABLE |                         |           |
| CYLINDER PRESSURE                       | 2000 PSIG            |                         |           |
| CERTIFICATION DATE                      | 08/08/05             |                         |           |
| EXPIRATION DATE                         | 08/08/08             | TERM                    | 36 MONTHS |

ANALYZED BY

ISMAEL RANGSIYAWONG

CERTIFIED BY

HELENA TRAN

#### IMPORTANT

Information contained herein has been prepared at your request by qualified experts within Praxair Distribution, Inc. While we believe that the information is accurate within the limits of the analytical methods employed and is complete to the extent of the specific analyses performed, we make no warranty or representation as to the suitability of the use of the information for any particular purpose. The information is offered with the understanding that any use of the information is at the sole discretion and risk of the user. In no event shall the liability of Praxair Distribution, Inc., arising out of the use of the information contained herein exceed the fee established for providing such information.

## ***Appendix C***

### ***Process Data***

- A-50
- A-51

# REDWOOD FLARE A-50 SOURCE TEST

| Date         | Time     | Flow<br>SCFM |     | Temp<br>°F |      |    |
|--------------|----------|--------------|-----|------------|------|----|
|              |          | MIN          | MAX | MIN        | MAX  |    |
| 2006/06/07   | 08:12:00 | 861          | 884 | 1638       | 1659 | R1 |
| 2006/06/07   | 08:14:00 | 859          | 886 | 1642       | 1652 | R1 |
| 2006/06/07   | 08:16:00 | 859          | 883 | 1640       | 1650 | R1 |
| 2006/06/07   | 08:18:00 | 863          | 889 | 1640       | 1655 | R1 |
| 2006/06/07   | 08:20:00 | 867          | 895 | 1641       | 1650 | R1 |
| 2006/06/07   | 08:22:00 | 859          | 884 | 1639       | 1651 | R1 |
| 2006/06/07   | 08:24:00 | 853          | 886 | 1638       | 1652 | R1 |
| 2006/06/07   | 08:26:00 | 864          | 888 | 1636       | 1657 | R1 |
| 2006/06/07   | 08:28:00 | 858          | 885 | 1643       | 1659 | R1 |
| 2006/06/07   | 08:30:00 | 869          | 884 | 1638       | 1657 | R1 |
| 2006/06/07   | 08:32:00 | 866          | 889 | 1640       | 1657 | R1 |
| 2006/06/07   | 08:34:00 | 856          | 885 | 1639       | 1657 | R1 |
| 2006/06/07   | 08:36:00 | 858          | 885 | 1641       | 1655 | R1 |
| 2006/06/07   | 08:38:00 | 859          | 892 | 1638       | 1652 | R1 |
| 2006/06/07   | 08:40:00 | 863          | 892 | 1641       | 1655 | R1 |
| Avg          |          | 861          | 887 | 1640       | 1655 |    |
| Combined Avg |          |              | 874 |            | 1647 |    |
| 2006/06/07   | 08:42:00 | 856          | 880 | 1637       | 1650 |    |
| 2006/06/07   | 08:44:00 | 860          | 884 | 1639       | 1653 |    |
| 2006/06/07   | 08:46:00 | 858          | 884 | 1643       | 1655 |    |
| 2006/06/07   | 08:48:00 | 854          | 885 | 1644       | 1650 |    |
| 2006/06/07   | 08:50:00 | 863          | 889 | 1639       | 1655 |    |
| 2006/06/07   | 08:52:00 | 857          | 883 | 1638       | 1656 | R2 |
| 2006/06/07   | 08:54:00 | 862          | 886 | 1642       | 1656 | R2 |
| 2006/06/07   | 08:56:00 | 861          | 896 | 1634       | 1654 | R2 |
| 2006/06/07   | 08:58:00 | 860          | 893 | 1642       | 1656 | R2 |
| 2006/06/07   | 09:00:00 | 856          | 882 | 1640       | 1651 | R2 |
| 2006/06/07   | 09:02:00 | 861          | 885 | 1639       | 1657 | R2 |
| 2006/06/07   | 09:04:00 | 860          | 891 | 1639       | 1655 | R2 |
| 2006/06/07   | 09:06:00 | 861          | 883 | 1643       | 1653 | R2 |
| 2006/06/07   | 09:08:00 | 855          | 881 | 1641       | 1652 | R2 |
| 2006/06/07   | 09:10:00 | 859          | 880 | 1640       | 1654 | R2 |
| 2006/06/07   | 09:12:00 | 862          | 889 | 1640       | 1655 | R2 |
| 2006/06/07   | 09:14:00 | 854          | 880 | 1642       | 1653 | R2 |
| 2006/06/07   | 09:16:00 | 863          | 882 | 1642       | 1651 | R2 |
| 2006/06/07   | 09:18:00 | 863          | 884 | 1642       | 1655 | R2 |
| 2006/06/07   | 09:20:00 | 860          | 887 | 1640       | 1658 | R2 |
| 2006/06/07   | 09:22:00 | 857          | 883 | 1641       | 1653 | R2 |
| Avg          |          | 859          | 885 | 1640       | 1654 |    |
| Combined Avg |          |              | 872 |            | 1647 |    |
| 2006/06/07   | 09:24:00 | 860          | 876 | 1644       | 1653 |    |
| 2006/06/07   | 09:26:00 | 858          | 887 | 1639       | 1653 |    |
| 2006/06/07   | 09:28:00 | 861          | 884 | 1640       | 1659 |    |
| 2006/06/07   | 09:30:00 | 858          | 883 | 1645       | 1655 |    |
| 2006/06/07   | 09:32:00 | 859          | 885 | 1640       | 1654 | R3 |
| 2006/06/07   | 09:34:00 | 860          | 885 | 1635       | 1653 | R3 |
| 2006/06/07   | 09:36:00 | 863          | 883 | 1640       | 1657 | R3 |
| 2006/06/07   | 09:38:00 | 862          | 881 | 1642       | 1658 | R3 |
| 2006/06/07   | 09:40:00 | 862          | 877 | 1642       | 1654 | R3 |
| 2006/06/07   | 09:42:00 | 852          | 886 | 1639       | 1654 | R3 |
| 2006/06/07   | 09:44:00 | 850          | 890 | 1639       | 1656 | R3 |
| 2006/06/07   | 09:46:00 | 856          | 883 | 1641       | 1657 | R3 |
| 2006/06/07   | 09:48:00 | 858          | 878 | 1641       | 1647 | R3 |
| 2006/06/07   | 09:50:00 | 853          | 875 | 1636       | 1658 | R3 |
| 2006/06/07   | 09:52:00 | 855          | 880 | 1636       | 1658 | R3 |
| 2006/06/07   | 09:54:00 | 853          | 883 | 1642       | 1651 | R3 |
| 2006/06/07   | 09:56:00 | 864          | 883 | 1645       | 1652 | R3 |
| 2006/06/07   | 09:58:00 | 864          | 883 | 1644       | 1657 | R3 |
| 2006/06/07   | 10:00:00 | 857          | 884 | 1642       | 1657 | R3 |
| Avg          |          | 858          | 882 | 1641       | 1655 |    |
| Combined Avg |          |              | 870 |            | 1648 |    |

## REDWOOD FLARE A-51 SOURCE TEST

| Date         | Time     | Temp<br>DEG. F |      | Flow<br>SCFM |      |    |
|--------------|----------|----------------|------|--------------|------|----|
|              |          | MIN            | MAX  | MIN          | MAX  |    |
| 2006/06/08   | 08:52:00 | 1421           | 1439 | 2596         | 2653 | R1 |
| 2006/06/08   | 08:54:00 | 1431           | 1445 | 2597         | 2644 | R1 |
| 2006/06/08   | 08:56:00 | 1445           | 1460 | 2601         | 2645 | R1 |
| 2006/06/08   | 08:58:00 | 1452           | 1467 | 2591         | 2637 | R1 |
| 2006/06/08   | 09:00:00 | 1446           | 1456 | 2592         | 2647 | R1 |
| 2006/06/08   | 09:02:00 | 1442           | 1450 | 2605         | 2641 | R1 |
| 2006/06/08   | 09:04:00 | 1440           | 1454 | 2610         | 2646 | R1 |
| 2006/06/08   | 09:06:00 | 1430           | 1454 | 2604         | 2636 | R1 |
| 2006/06/08   | 09:08:00 | 1430           | 1457 | 2609         | 2636 | R1 |
| 2006/06/08   | 09:10:00 | 1440           | 1458 | 2601         | 2640 | R1 |
| 2006/06/08   | 09:12:00 | 1453           | 1462 | 2599         | 2633 | R1 |
| 2006/06/08   | 09:14:00 | 1439           | 1455 | 2592         | 2622 | R1 |
| 2006/06/08   | 09:16:00 | 1430           | 1444 | 2597         | 2622 | R1 |
| 2006/06/08   | 09:18:00 | 1431           | 1453 | 2601         | 2635 | R1 |
| 2006/06/08   | 09:20:00 | 1453           | 1461 | 2599         | 2635 | R1 |
| 2006/06/08   | 09:22:00 | 1461           | 1469 | 2577         | 2633 | R1 |
| Avg          |          | 1440           | 1455 | 2598         | 2638 |    |
| Combined Avg |          |                | 1448 |              | 2618 |    |
| 2006/06/08   | 09:24:00 | 1468           | 1495 | 2597         | 2632 |    |
| 2006/06/08   | 09:26:00 | 1438           | 1492 | 2593         | 2629 |    |
| 2006/06/08   | 09:28:00 | 1431           | 1439 | 2601         | 2628 |    |
| 2006/06/08   | 09:30:00 | 1432           | 1489 | 2584         | 2640 |    |
| 2006/06/08   | 09:32:00 | 1433           | 1465 | 2599         | 2635 |    |
| 2006/06/08   | 09:34:00 | 1431           | 1437 | 2590         | 2633 | R2 |
| 2006/06/08   | 09:36:00 | 1435           | 1439 | 2587         | 2627 | R2 |
| 2006/06/08   | 09:38:00 | 1426           | 1439 | 2587         | 2635 | R2 |
| 2006/06/08   | 09:40:00 | 1424           | 1454 | 2580         | 2621 | R2 |
| 2006/06/08   | 09:42:00 | 1454           | 1468 | 2572         | 2613 | R2 |
| 2006/06/08   | 09:44:00 | 1460           | 1471 | 2596         | 2620 | R2 |
| 2006/06/08   | 09:46:00 | 1469           | 1477 | 2573         | 2620 | R2 |
| 2006/06/08   | 09:48:00 | 1453           | 1469 | 2580         | 2610 | R2 |
| 2006/06/08   | 09:50:00 | 1424           | 1453 | 2574         | 2616 | R2 |
| 2006/06/08   | 09:52:00 | 1424           | 1435 | 2576         | 2605 | R2 |
| 2006/06/08   | 09:54:00 | 1416           | 1426 | 2572         | 2606 | R2 |
| 2006/06/08   | 09:56:00 | 1422           | 1476 | 2571         | 2614 | R2 |
| 2006/06/08   | 09:58:00 | 1465           | 1479 | 2547         | 2610 | R2 |
| 2006/06/08   | 10:00:00 | 1464           | 1489 | 2539         | 2625 | R2 |
| 2006/06/08   | 10:02:00 | 1427           | 1482 | 2584         | 2606 | R2 |
| Avg          |          | 1440           | 1464 | 2580         | 2621 |    |
| Combined Avg |          |                | 1452 |              | 2601 |    |
| 2006/06/08   | 10:04:00 | 1423           | 1430 | 2575         | 2613 |    |
| 2006/06/08   | 10:06:00 | 1420           | 1426 | 2582         | 2609 |    |
| 2006/06/08   | 10:08:00 | 1418           | 1455 | 2587         | 2616 |    |
| 2006/06/08   | 10:10:00 | 1455           | 1472 | 2587         | 2616 |    |
| 2006/06/08   | 10:12:00 | 1468           | 1474 | 2574         | 2604 |    |
| 2006/06/08   | 10:14:00 | 1455           | 1476 | 2581         | 2616 | R3 |
| 2006/06/08   | 10:16:00 | 1440           | 1455 | 2581         | 2605 | R3 |
| 2006/06/08   | 10:18:00 | 1440           | 1446 | 2583         | 2608 | R3 |
| 2006/06/08   | 10:20:00 | 1442           | 1560 | 2587         | 2615 | R3 |
| 2006/06/08   | 10:22:00 | 1429           | 1527 | 2572         | 2608 | R3 |
| 2006/06/08   | 10:24:00 | 1442           | 1469 | 2585         | 2619 | R3 |
| 2006/06/08   | 10:26:00 | 1420           | 1459 | 2569         | 2612 | R3 |
| 2006/06/08   | 10:28:00 | 1373           | 1469 | 2571         | 2606 | R3 |
| 2006/06/08   | 10:30:00 | 1369           | 1430 | 2547         | 2603 | R3 |
| 2006/06/08   | 10:32:00 | 1408           | 1426 | 2561         | 2597 | R3 |
| 2006/06/08   | 10:34:00 | 1419           | 1428 | 2580         | 2608 | R3 |
| 2006/06/08   | 10:36:00 | 1419           | 1433 | 2572         | 2607 | R3 |
| 2006/06/08   | 10:38:00 | 1433           | 1457 | 2566         | 2595 | R3 |
| 2006/06/08   | 10:40:00 | 1457           | 1476 | 2569         | 2610 | R3 |
| 2006/06/08   | 10:42:00 | 1476           | 1558 | 2570         | 2597 | R3 |
| 2006/06/08   | 10:44:00 | 1453           | 1553 | 2569         | 2603 | R3 |
| Avg          |          | 1431           | 1470 | 2575         | 2608 |    |
| Combined Avg |          |                | 1451 |              | 2591 |    |

## Landfill Process Data Sheet

Landfill: Redwood

Test date: 6-7-06

Source: Flare, IC Engine, Turbine  
circle one

| Landfill Gas Data         | Run 1 |      | Run 2  |      | Run 3 |  |
|---------------------------|-------|------|--------|------|-------|--|
| Time                      | 811   | 2841 | 2852   | 0922 | 0131  |  |
| Methane (%)               | 51.6  |      |        |      |       |  |
| HHV based on methane      |       |      |        |      |       |  |
| F-factor based on methane |       |      |        |      |       |  |
| O2 (%)                    | 1.6   |      |        |      |       |  |
| Landfill Gas Rate (scfm)  | 870   |      | 865879 |      | 868   |  |
| Engine (KW)               |       |      |        |      |       |  |
| Condensate Rate (gpm)     |       |      |        |      |       |  |
| Flare Temp (F)            | 1650  |      | 1647   |      | 1649  |  |

Thermocouple Measurement Location Top Middle Bottom  
circle one

HHV = methane (%) x 1018

Set point 1650

# Landfill Process Data Sheet

Landfill: Redwood Landf. II A 51

Test date: 06-08-06

Source: Flare, IC Engine, Turbine

circle one

| Landfill Gas Data         | Run 1 |  | Run 2 |  | Run 3 |  |
|---------------------------|-------|--|-------|--|-------|--|
| Time                      | 0852  |  | 0933  |  |       |  |
| Methane (%)               | 48.6  |  |       |  |       |  |
| HHV based on methane      |       |  |       |  |       |  |
| F-factor based on methane |       |  |       |  |       |  |
| O2 (%)                    | 1.9   |  |       |  |       |  |
| Landfill Gas Rate (scfm)  | 2620  |  | 2608  |  | 2580  |  |
| Engine (KW)               |       |  |       |  |       |  |
| Condensate Rate (gpm)     |       |  |       |  |       |  |
| Flare Temp (F)            | 1450  |  | 1468  |  | 1429  |  |

Thermocouple Measurement Location

circle one

Top

Middle

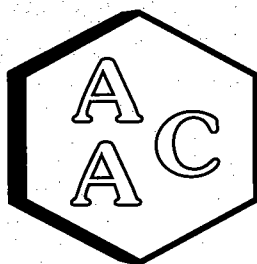
Bottom

HHV = methane (%) x 1018

See pair 1450

## ***Appendix D***

### ***Analytical Results***



## Atmospheric Analysis & Consulting, Inc.

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CLIENT : Shaw Environmental, Inc.  
PROJECT NAME : Redwood Landfill  
AAC PROJECT NO. : 060263  
REPORT DATE : 06/20/2006

On June 9, 2006, Atmospheric Analysis & Consulting, Inc. received one (1) Tedlar Bag for BTU Analysis, one (1) Tedlar Bag for BTU Analysis and Volatile Organic Compounds Analysis by EPA Method 25C, and one (1) Tedlar Bag to be put on hold. Upon receipt the samples were assigned unique Laboratory ID numbers as follows:

| Client ID | Lab No.      |
|-----------|--------------|
| A50, R2   | 060263-15746 |
| A51, R2   | 060263-15747 |
| A51 Dup   | 060263-15748 |

BTU Analysis includes the following tests: Sulfur Analysis by ASTM Method D-5504, C1-C6+ Analysis by EPA Method 18, and Fixed Gases Analysis by EPA Method 3C.

EPA 25C Analysis - Up to a 1 ml aliquot of gaseous sample is injected into the GC/FID for analysis following EPA25C as specified in the SOW.

ASTM D-5504 - Up to a 1 ml aliquot of sample is injected into the GC/SCD for analysis following ASTM D-5504 as specified in the SOW.

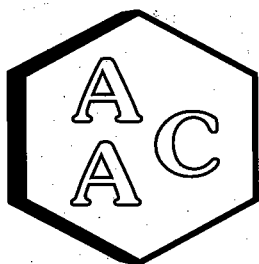
EPA 18 Analysis - Up to a 1 ml aliquot of samples is injected into the GC/FID for analysis following EPA 18 as specified in the SOW.

EPA 3C Analysis - Up to a 1 ml aliquot of samples is injected into the GC/TCD for analysis following EPA 3C as specified in the SOW.

No problems were encountered during receiving, preparation, and/ or analysis of these samples. The test results included in this report meet all requirements of the NELAC Standards and/or AAC SOP# AACI- EPA 25C, ASTM D-5504, EPA 18 and EPA 3C.

I certify that this data is technically accurate, complete, and in compliance with the terms and conditions of the contract. Release of the data contained in this hardcopy data package and its electronic data deliverable submitted on diskette has been authorized by the Laboratory Director or his designee, as verified by the following signature.

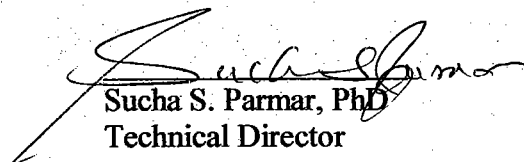




## Atmospheric Analysis & Consulting, Inc.

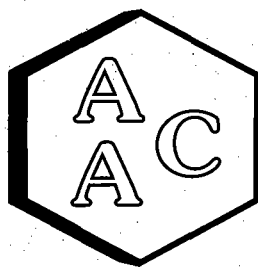
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If you have any question or require further explanation of data results, please contact the undersigned.

  
Sucha S. Parmar, PhD  
Technical Director

This report consists of 11 pages.





# Atmospheric Analysis & Consulting, Inc.

Page 3

## Laboratory Analysis Report

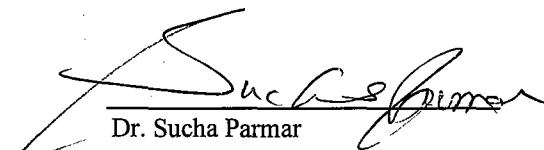
Client: : Shaw Environmental, Inc.  
Project No. : 060263  
Matrix : air  
Units : %

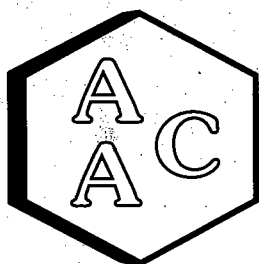
Sampling Date : 06/7-8/06  
Receiving Date : 06/09/2006  
Analysis Date : 06/09/2006  
Report Date : 06/20/2006

### EPA Method 3C

| Detection Limit: 0.1 % |              | Analyte  |        |          |    |         |      |
|------------------------|--------------|----------|--------|----------|----|---------|------|
| Client ID              | AAC ID       | Hydrogen | Oxygen | Nitrogen | CO | Methane | CO2  |
| A50, R2                | 060263-15746 | ND       | 2.5    | 12.6     | ND | 51.3    | 33.6 |
| A51, R2                | 060263-15747 | ND       | 3.2    | 15.4     | ND | 48.7    | 32.7 |

\* Per request on the COC, the duplicate sample for A51,R2 was analyzed because the O2 was above 3%.  
The duplicate sample also had an O2 value of 3.4 %.

  
Dr. Sucha Parmar  
Technical Director



# Atmospheric Analysis & Consulting, Inc.

## Laboratory Analysis Report

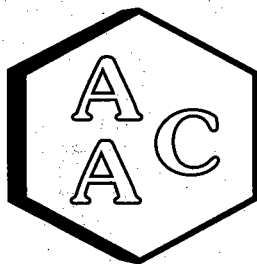
CLIENT : Shaw Environmental, Inc.  
PROJECT NO. : 060263  
MATRIX : GAS

RECEIVING DATE : 06/09/2006  
ANALYSIS DATE : 06/09/2006  
REPORT DATE : 06/20/2006

|                                 | Client ID:       | A50, R2      | A51, R2      | Detection Limits |
|---------------------------------|------------------|--------------|--------------|------------------|
|                                 | AAC ID:          | 060263-15746 | 060263-15747 |                  |
|                                 | Compounds, Units |              |              |                  |
| H <sub>2</sub> S, ppmv          |                  | 92           | 326          | 0.5ppmv          |
| Carbonyl Sulfide, ppmv          |                  | ND           | ND           | 0.5ppmv          |
| Methyl Mercaptan, ppmv          |                  | ND           | ND           | 0.5ppmv          |
| Ethyl Mercaptan, ppmv           |                  | ND           | ND           | 0.5ppmv          |
| Propyl Mercaptan, ppmv          |                  | ND           | ND           | 0.5ppmv          |
| Carbon Disulfide, ppmv          |                  | ND           | ND           | 0.5ppmv          |
| Dimethyl sulfide, ppmv          |                  | ND           | ND           | 0.5ppmv          |
| Allyl Sulfide, ppmv             |                  | ND           | ND           | 0.5ppmv          |
| Propyl Sulfide, ppmv            |                  | ND           | ND           | 0.5ppmv          |
| Allyl Disulfide, ppmv           |                  | ND           | ND           | 0.5ppmv          |
| Butyl Sulfide, ppmv             |                  | ND           | ND           | 0.5ppmv          |
| Ethyl Methyl Sulfide, ppmv      |                  | ND           | ND           | 0.5ppmv          |
| Thiophene, ppmv                 |                  | ND           | ND           | 0.5ppmv          |
| Dimethyl Disulfide, ppmv        |                  | ND           | ND           | 0.5ppmv          |
| Butyl Mercaptan, ppmv           |                  | ND           | ND           | 0.5ppmv          |
| Allyl mercaptan, ppmv           |                  | ND           | ND           | 0.5ppmv          |
| C1 hydrocarbons, ppmv           |                  | 511730       | 488215       | 0.3 ppmv         |
| C2 hydrocarbons, ppmv           |                  | <30          | <30          | 0.3 ppmv         |
| C3 hydrocarbons, ppmv           |                  | 9.4          | 10           | 0.3 ppmv         |
| C4 hydrocarbons, ppmv           |                  | 7.1          | 6.8          | 0.3 ppmv         |
| C5 hydrocarbons, ppmv           |                  | 3.6          | 2.2          | 0.3 ppmv         |
| C6 hydrocarbons, ppmv           |                  | 6.8          | 2.6          | 0.3 ppmv         |
| C6 + hydrocarbons, ppmv         |                  | 1.2          | 1.8          | 0.3 ppmv         |
| CO <sub>2</sub> , %             |                  | 34           | 33           | 0.1 %            |
| CO, %                           |                  | ND           | ND           | 0.1 %            |
| O <sub>2</sub> , %              |                  | 2.5          | 3.2          | 0.1 %            |
| N <sub>2</sub> , %              |                  | 13           | 15           | 0.1 %            |
| H <sub>2</sub> , %              |                  | ND           | ND           | 0.1 %            |
| F Factor(dscf Exhaust/MM Btu)   |                  | 9147         | 9158         |                  |
| Total Wt.% Adjusted Sp. Gravity |                  | 0.9          | 0.9539       |                  |
| FUEL GAS BTU per LBM            |                  | 12221        | 11660.9      |                  |
| FUEL GAS BTU per CU. FT         |                  | 519          | 495          |                  |

Dr. Sucha Parmar  
Technical Director





## Atmospheric Analysis & Consulting, Inc.

### LABORATORY ANALYSIS REPORT

CLIENT : Shaw Environmental, Inc.  
PROJECT NO. : 060263  
UNITS : PPMV

Sampling Date : 06/7-8/06  
Receiving Date : 06/09/2006  
Analysis Date : 06/09/2006  
Report Date : 06/20/2006

#### Total Reduced Sulfur Compounds Analysis by ASTM D-5504

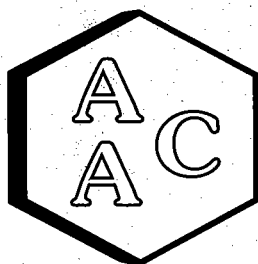
| Client ID.                                     | A50, R2      | A51, R2      | MDL  |
|------------------------------------------------|--------------|--------------|------|
| AAC ID                                         | 060263-15746 | 060263-15747 |      |
| Analysis Dilution Factor                       | 1.0 & 20     | 1.0 & 20     |      |
| Can Dilution Factor                            | 1.0          | 1.0          |      |
| H <sub>2</sub> S                               | 91           | 327          | 0.50 |
| Carbonyl Sulfide                               | <PQL         | <PQL         | 0.50 |
| SO <sub>2</sub>                                | <PQL         | <PQL         | 0.50 |
| Methyl Mercaptan                               | <PQL         | <PQL         | 0.50 |
| Ethyl Mercaptan                                | <PQL         | <PQL         | 0.50 |
| Dimethyl Sulfide                               | <PQL         | <PQL         | 0.50 |
| n-Butyl mercaptan                              | <PQL         | <PQL         | 0.50 |
| Carbon Disulfide                               | <PQL         | <PQL         | 0.50 |
| Allyl Sulfide                                  | <PQL         | <PQL         | 0.50 |
| Propyl Sulfide                                 | <PQL         | <PQL         | 0.50 |
| Allyl disulfide                                | <PQL         | <PQL         | 0.50 |
| Isopropyl Mercaptan                            | <PQL         | 1.8          | 0.50 |
| t-Butyl mercaptan                              | <PQL         | <PQL         | 0.50 |
| Propyl Mercaptan                               | <PQL         | <PQL         | 0.50 |
| Butyl Sulfide                                  | <PQL         | <PQL         | 0.50 |
| Ethyl methyl sulfide                           | <PQL         | <PQL         | 0.50 |
| Thiophene                                      | <PQL         | <PQL         | 0.50 |
| Isobutyl mercaptan                             | <PQL         | <PQL         | 0.50 |
| Dimethyl disulfide                             | <PQL         | <PQL         | 0.50 |
| Allyl mercaptan                                | <PQL         | <PQL         | 0.50 |
| 3-Methylthiophene                              | <PQL         | <PQL         | 0.50 |
| Tetrahydrothiophene                            | <PQL         | <PQL         | 0.50 |
| Diethyl sulfide                                | <PQL         | <PQL         | 0.50 |
| 2-Ethylthiophene                               | <PQL         | <PQL         | 0.50 |
| 2,5-Dimethylthiophene                          | <PQL         | <PQL         | 0.50 |
| Diethyl disulfide                              | <PQL         | <PQL         | 0.50 |
| Total Unidentified Sulfurs as H <sub>2</sub> S | <PQL         | 1.8          | 0.50 |
| Total Reduced Sulfurs                          | 92           | 331          | 0.50 |

PQL = Practical Quantitation Limit (MDL x Analysis Dilution factor)

All compounds concentrations expressed in terms of H<sub>2</sub>S.

  
Dr. Sucha Parmar  
Technical Director





## Atmospheric Analysis & Consulting, Inc.

### Laboratory Analysis Report

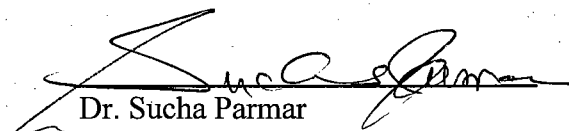
Client : Shaw Environmental, Inc.  
Project No. : 060263  
Matrix : air  
Units : ppmv

Sampling Date : 06/07/2006  
Receiving Date : 06/09/2006  
Analysis Date : 06/09/2006  
Report Date : 06/20/2006

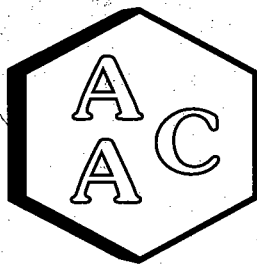
#### EPA Method 25C

| Detection Limit: |              | 0.3 ppmv |
|------------------|--------------|----------|
| Client Sample ID | AAC ID       | NMHC**   |
| A50, R2          | 060263-15746 | 456      |
| A51, R2          | 060263-15747 | 213      |

\*\*Non-Methane Hydrocarbons as methane

  
Dr. Sucha Parmar  
Technical Director





## Atmospheric Analysis & Consulting, Inc.

### Quality Control/Quality Assurance Report

DATE ANALYZED : 06/09/2006  
UNITS : ppmv

#### I - Method Blank

| AAC ID       | Analyte | MB Conc. |
|--------------|---------|----------|
| Method Blank | H2S     | ND       |

#### II-Laboratory Control Spike - ASTM D-5504

| Analyte | Spike Added | Spike Res | Dup Spike Res | Spike % Rec * | Spike Dup % Rec * | RPD** |
|---------|-------------|-----------|---------------|---------------|-------------------|-------|
| H2S     | 5.00        | 4.95      | 5.09          | 99            | 102               | 2.8   |

#### III - Duplicate Analysis - ASTM D-5504 - 060263-15746 (20x)

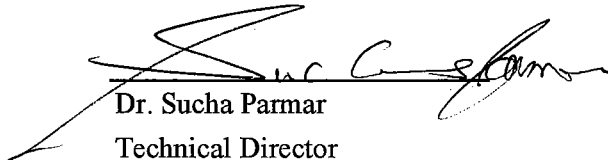
| Analyte | Sample Conc | Duplicate Conc | Mean | % RPD |
|---------|-------------|----------------|------|-------|
| H2S     | 99.0        | 83.4           | 91.2 | 17    |

#### IV - Matrix Spike - ASTM D-5504 - 060256-15710 (40x)

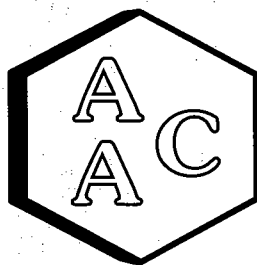
| Analyte | Sample Conc. | Spike Added | Spike Res | Dup Spike Res | Spike % Rec * | Spike Dup % Rec * | RPD** |
|---------|--------------|-------------|-----------|---------------|---------------|-------------------|-------|
| H2S     | 2.28         | 2.50        | 4.66      | 4.19          | 95            | 76                | 22    |

\* Must be 70-130%

\*\* Must be  $\leq 25\%$

  
Dr. Sucha Parmar  
Technical Director





# Atmospheric Analysis & Consulting, Inc.

## Quality Control/Quality Assurance Report

Date Analyzed : 06/09/2006

EPA Method 25.3

### I - Method Blank

| Analyte | Units | Sample Analysis |
|---------|-------|-----------------|
| NMHC    | ppm   | ND              |

### II - Duplicate Analysis

| AAC ID                | Analyte | Units | Sample Analysis | Duplicate Analysis | Mean  | % RPD ** |
|-----------------------|---------|-------|-----------------|--------------------|-------|----------|
| 060262-15741<br>(10x) | NMHC    | ppm   | 481.3           | 481.1              | 481.2 | 0.0      |

### III - Matrix Spikes

| AAC ID                | Analyte | Units | Sample Conc. | Spike Added | Sample Spike Conc. | Sample Spike Dup. Conc. | Spike % Rec.* | Spike Dup % Rec.* | % RPD ** |
|-----------------------|---------|-------|--------------|-------------|--------------------|-------------------------|---------------|-------------------|----------|
| 060262-15741<br>(10x) | NMHC    | ppm   | 240.6        | 125         | 374                | 391                     | 107           | 120               | 4.5      |

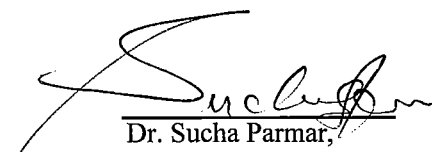
### IV - 1 ppm Backflush

| Analyte | Units | Sample Analysis | Sample Analysis dup | Sample % Rec.*** | Sample Dup % Rec. *** |
|---------|-------|-----------------|---------------------|------------------|-----------------------|
| NMHC    | ppm   | 1.05            | 1.05                | 105              | 105                   |

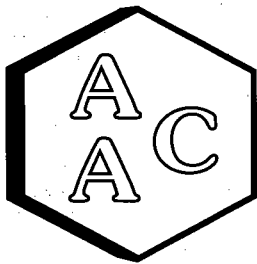
\* Must be 90-110%

\*\* Must be 75-125%

\*\*\* Must be 80-120%

  
Dr. Sucha Parmar,  
Technical Director





# Atmospheric Analysis & Consulting, Inc.

## Quality Control/Quality Assurance Report

Date Analyzed : 06/09/2006

### I-Method Blank - EPA 3C

|              | Analyte  | Units | Sample Analysis |
|--------------|----------|-------|-----------------|
| Method Blank | Hydrogen | %     | ND              |
|              | Oxygen   | %     | ND              |
|              | Nitrogen | %     | ND              |
|              | CO       | %     | ND              |
|              | Methane  | %     | ND              |
|              | CO2      | %     | ND              |

### II-LCS - EPA 3C

|     | Analyte  | Spike Conc. | LCS Conc. | LCS Dup Conc. | % Rec | % RPD |
|-----|----------|-------------|-----------|---------------|-------|-------|
| LCS | Hydrogen | 20          | 19        | 19            | 95    | 3.9   |
|     | Nitrogen | 20          | 23        | 23            | 116   | 0.3   |
|     | CO       | 20          | 20        | 20            | 101   | 1.7   |
|     | Methane  | 20          | 23        | 23            | 114   | 1.2   |
|     | CO2      | 20          | 22        | 21            | 108   | 1.7   |

### III-Duplicate Analysis -EPA 3C

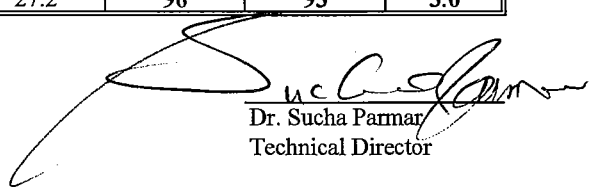
| AAC ID       | Analyte  | Sample Analysis | Duplicate Analysis | Mean | % RPD |
|--------------|----------|-----------------|--------------------|------|-------|
| 060263-15746 | Hydrogen | 0.0             | 0.0                | 0.0  | NA    |
|              | Oxygen   | 2.7             | 2.7                | 2.7  | 0.6   |
|              | Nitrogen | 13.3            | 13.5               | 13.4 | 0.6   |
|              | CO       | 0.0             | 0.0                | 0.0  | NA    |
|              | Methane  | 54.3            | 55.1               | 54.7 | 0.7   |
|              | CO2      | 35.6            | 35.9               | 35.8 | 0.4   |

### IV-Matrix Spike Analysis -EPA 3C

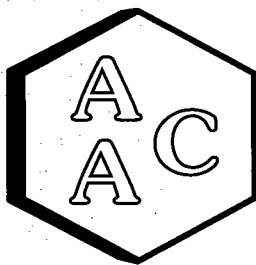
| AAC ID            | Analyte  | Sample Conc. | Spike Added | Sample Spike Conc. | Sample Spike Dupl. Conc. | Spike % Rec. * | Spike Dupl % Rec. * | % RPD ** |
|-------------------|----------|--------------|-------------|--------------------|--------------------------|----------------|---------------------|----------|
| 060263-15746 (2x) | Hydrogen | 0.0          | 10.0        | 8.7                | 8.5                      | 87             | 85                  | 3.3      |
|                   | Nitrogen | 6.7          | 10.0        | 17.4               | 17.6                     | 107            | 109                 | 1.4      |
|                   | CO       | 0.0          | 10.0        | 10.1               | 10.0                     | 101            | 100                 | 1.3      |
|                   | Methane  | 27.3         | 10.0        | 37.3               | 36.8                     | 99             | 95                  | 4.7      |
|                   | CO2      | 17.9         | 10.0        | 27.5               | 27.2                     | 96             | 93                  | 3.0      |

\* Must be 70-130%

\*\* Must be  $\leq 25\%$

  
Dr. Sucha Parmar  
Technical Director





# Atmospheric Analysis & Consulting, Inc.

## Quality Control/Quality Assurance Report

Date Analyzed : 06/09/2006  
EPA Method 18

### I - Method Blank

| AAC ID       | Analyte | Units | MB Conc. |
|--------------|---------|-------|----------|
| Method Blank | Methane | ppmv  | ND       |
|              | Ethane  | ppmv  | ND       |
|              | Propane | ppmv  | ND       |
|              | Butane  | ppmv  | ND       |
|              | Pentane | ppmv  | ND       |
|              | Hexane  | ppmv  | ND       |

### II - LCS

| AAC ID               | Analyte | Theoretical Conc. | LCS Conc. | LCS Conc. Dup | % Rec | % RPD |
|----------------------|---------|-------------------|-----------|---------------|-------|-------|
| Lab Control Standard | Methane | 100.2             | 103.7     | 106.9         | 105   | 3.1   |
|                      | Ethane  | 100.6             | 102.4     | 106.6         | 104   | 4.0   |
|                      | Propane | 100.0             | 102.4     | 108.0         | 105   | 5.3   |
|                      | Butane  | 99.6              | 99.8      | 105.5         | 103   | 5.6   |
|                      | Pentane | 99.6              | 96.7      | 102.9         | 100   | 6.2   |
|                      | Hexane  | 97.6              | 90.0      | 99.5          | 97    | 10    |

### III - Duplicate Analysis

| AAC ID                  | Analyte | Units | Sample Conc. | Duplicate Conc. | Mean   | %RPD |
|-------------------------|---------|-------|--------------|-----------------|--------|------|
| 060262-15743<br>(2000x) | Methane | ppmv  | 444503       | 451864          | 448200 | 1.6  |
|                         | Ethane  | ppmv  | 0.0          | 0.0             | 0.0    | 0.0  |
|                         | Propane | ppmv  | 0.0          | 0.0             | 0.0    | 0.0  |
|                         | Butane  | ppmv  | 0.0          | 0.0             | 0.0    | 0.0  |
|                         | Pentane | ppmv  | 0.0          | 0.0             | 0.0    | 0.0  |
|                         | Hexane  | ppmv  | 0.0          | 0.0             | 0.0    | 0.0  |

### IV - Spiked Sample

| AAC ID                  | Analyte | Units | Sample Conc. | Spike Added | Spike Res | Dup Spike Res | Spike % Rec * | Spike Dup % Rec * | RPD** % |
|-------------------------|---------|-------|--------------|-------------|-----------|---------------|---------------|-------------------|---------|
| 060262-15743<br>(4000x) | Methane | ppmv  | 112          | 100.0       | 239       | 243           | 127           | 131 (M7)          | 3.1     |
|                         | Ethane  | ppmv  | 0.0          | 100.0       | 119.6     | 120.5         | 120           | 121               | 0.7     |
|                         | Propane | ppmv  | 0.0          | 100.0       | 121.2     | 121.1         | 121           | 121               | 0.1     |
|                         | Butane  | ppmv  | 0.0          | 100.0       | 119.2     | 119.4         | 119           | 119               | 0.2     |
|                         | Pentane | ppmv  | 0.0          | 100.0       | 118.2     | 117.3         | 118           | 117               | 0.8     |
|                         | Hexane  | ppmv  | 0.0          | 100.0       | 115.7     | 111.9         | 116           | 112               | 3.4     |

\* Must be 70-130%

\*\* Must be ≤ 25%

M7 - Matrix spike recovery was above laboratory acceptance limits, the MB, LCS, LCSD had good recoveries and RPD's, therefore high recovery due to matrix effect.

Dr. Sucha Parmar  
Technical Director



AAC

060263

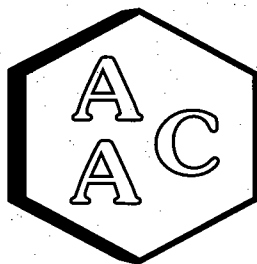
**CLIENT**

Redwood Landfill

DATE: 6-8-06

PAGE: 1 of 1

[illegible]



## Atmospheric Analysis & Consulting, Inc.

CLIENT : Shaw Environmental, Inc.  
PROJECT NAME : Redwood Landfill  
AAC PROJECT NO. : 060263  
REPORT DATE : 06/20/06

On June 9, 2006, Atmospheric Analysis & Consulting, Inc. received one (1) Tedlar bag for Volatile Organic Compounds analysis by EPA method TO-15. Upon receipt the sample was assigned a unique Laboratory ID number as follows:

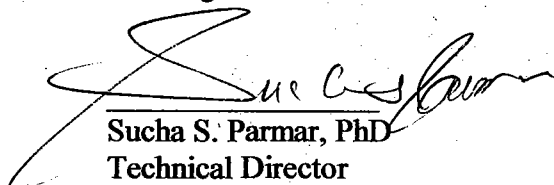
| Client ID | Lab ID       |
|-----------|--------------|
| A51, R2   | 060263-15747 |

TO-14/15 Analysis - Up to a 500 ml aliquot of samples is concentrated, put through a water and CO2 management system, cryofocused and injected into the GC/MS (full scan mode) for analysis following EPA Method TO-14/15 as specified in the SOW.

No problems were encountered during receiving, preparation and/ or analysis of these samples. The test results included in this report meet all requirements of the NELAC Standards and/or AAC SOP# AACI-TO-15/PAMS. Estimated uncertainty of the test results will be provided upon request.

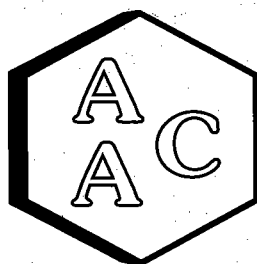
I certify that this data is technically accurate, complete and in compliance with the terms and conditions of the contract. The Laboratory Director or his designee, as verified by the following signature, has authorized the release of the data contained in this hardcopy data package.

If you have any questions or require further explanation of data results, please contact the undersigned.

  
Sucha S. Parmar, PhD  
Technical Director

This report consists of 22 pages.





# Atmospheric Analysis & Consulting, Inc.

## Laboratory Analysis Report

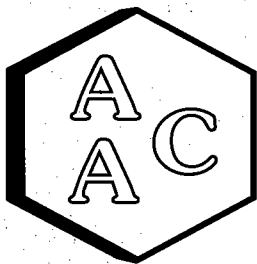
CLIENT : Shaw Environmental  
PROJECT NO : 060263  
MATRIX : AIR  
UNITS : PPB (v/v)

DATE RECEIVED : 06/09/06  
DATE REPORTED : 06/20/06

### VOLATILE ORGANIC COMPOUNDS BY EPA TO-15

| Client ID                              | A51, R2      |           |           | Sample<br>Reporting<br>Limit<br>(RLxDF's) | Method<br>Reporting<br>Limit |
|----------------------------------------|--------------|-----------|-----------|-------------------------------------------|------------------------------|
| AAC ID                                 | 060263-15747 |           |           |                                           |                              |
| Date Sampled                           | 6/7/2006     |           |           |                                           |                              |
| Date Analyzed                          | 6/17/2006    |           |           |                                           |                              |
| Can Dilution Factor                    | 1.00         |           |           |                                           |                              |
|                                        | Result       | Qualifier | Dil. Fac. |                                           |                              |
| Chlorodifluoromethane                  | 575          |           | 100       | 100                                       | 1.0                          |
| Propylene                              | 7960         |           | 1000      | 1000                                      | 1.0                          |
| Dichlorodifluoromethane                | 642          |           | 100       | 100                                       | 1.0                          |
| Chloromethane                          | ND           | U         | 100       | 100                                       | 1.0                          |
| 1,2-Dichloro-1,1,2,2-Tetrafluoroethane | ND           | U         | 100       | 100                                       | 1.0                          |
| Vinyl Chloride                         | 431          |           | 100       | 100                                       | 1.0                          |
| Methanol                               | ND           | U         | 100       | 500                                       | 5.0                          |
| 1,3-Butadiene                          | ND           | U         | 100       | 100                                       | 1.0                          |
| Bromomethane                           | ND           | U         | 100       | 100                                       | 1.0                          |
| Chloroethane                           | ND           | U         | 100       | 100                                       | 1.0                          |
| Dichlorofluoromethane                  | ND           | U         | 100       | 100                                       | 1.0                          |
| Ethanol                                | 294          |           | 100       | 200                                       | 2.0                          |
| Vinyl Bromide                          | ND           | U         | 100       | 100                                       | 1.0                          |
| Acetone                                | 1070         |           | 100       | 200                                       | 2.0                          |
| Trichlorofluoromethane                 | ND           | U         | 100       | 100                                       | 1.0                          |
| Isopropyl Alcohol                      | 269          |           | 100       | 200                                       | 2.0                          |
| Acrylonitrile                          | ND           | U         | 100       | 100                                       | 1.0                          |
| 1,1-Dichloroethylene                   | ND           | U         | 100       | 100                                       | 1.0                          |
| Methylene Chloride                     | ND           | U         | 100       | 100                                       | 1.0                          |
| Allyl Chloride (Chloroprene)           | ND           | U         | 100       | 100                                       | 1.0                          |
| Carbon Disulfide                       | ND           | U         | 100       | 100                                       | 1.0                          |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane  | ND           | U         | 100       | 100                                       | 1.0                          |
| t-1,2-Dichloroethylene                 | ND           | U         | 100       | 100                                       | 1.0                          |
| 1,1-Dichloroethane                     | ND           | U         | 100       | 100                                       | 1.0                          |
| MTBE                                   | ND           | U         | 100       | 100                                       | 1.0                          |
| Vinyl Acetate                          | ND           | U         | 100       | 100                                       | 1.0                          |
| 2-Butanone (MEK)                       | 1200         |           | 100       | 100                                       | 1.0                          |
| cis-1,2- Dichloroethene                | 174          |           | 100       | 100                                       | 1.0                          |
| Hexane                                 | 471          |           | 100       | 100                                       | 1.0                          |
| Chloroform                             | ND           | U         | 100       | 100                                       | 1.0                          |
| Ethyl Acetate                          | ND           | U         | 100       | 100                                       | 1.0                          |
| Tetrahydrofuran                        | 1500         |           | 100       | 100                                       | 1.0                          |
| 1,2-Dichloroethane                     | ND           | U         | 100       | 100                                       | 1.0                          |
| 1,1,1-Trichloroethane                  | ND           | U         | 100       | 100                                       | 1.0                          |





# Atmospheric Analysis & Consulting, Inc.

## Laboratory Analysis Report

CLIENT : Shaw Environmental  
PROJECT NO : 060263  
MATRIX : AIR  
UNITS : PPB (v/v)

DATE RECEIVED : 06/09/06  
DATE REPORTED : 06/20/06

### VOLATILE ORGANIC COMPOUNDS BY EPA TO-15

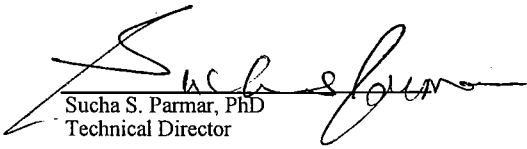
| Client ID                     | A51, R2      |           |           | Sample<br>Reporting<br>Limit<br>(RLxDf's) | Method<br>Reporting<br>Limit |
|-------------------------------|--------------|-----------|-----------|-------------------------------------------|------------------------------|
| AAC ID                        | 060263-15747 |           |           |                                           |                              |
| Date Sampled                  | 6/7/2006     |           |           |                                           |                              |
| Date Analyzed                 | 6/17/2006    |           |           |                                           |                              |
| Can Dilution Factor           | 1.00         |           |           |                                           |                              |
|                               | Result       | Qualifier | Dil. Fac. |                                           |                              |
| Benzene                       | 328          |           | 100       | 100                                       | 1.0                          |
| Carbon Tetrachloride          | ND           | U         | 100       | 100                                       | 1.0                          |
| Cyclohexane                   | 350          |           | 100       | 100                                       | 1.0                          |
| 1,2-Dichloropropane           | ND           | U         | 100       | 100                                       | 1.0                          |
| Bromodichloromethane          | ND           | U         | 100       | 100                                       | 1.0                          |
| 1,4-Dioxane                   | ND           | U         | 100       | 100                                       | 1.0                          |
| Trichloroethene               | ND           | U         | 100       | 100                                       | 1.0                          |
| 2,2,4-Trimethylpentane        | 168          |           | 100       | 100                                       | 1.0                          |
| Heptane                       | 830          |           | 100       | 100                                       | 1.0                          |
| cis-1,3-Dichloropropene       | ND           | U         | 100       | 100                                       | 1.0                          |
| 4-Methyl-2-Pentanone (MiBK)   | 153          |           | 100       | 100                                       | 1.0                          |
| t-1,3-Dichloropropene         | ND           | U         | 100       | 100                                       | 1.0                          |
| 1,1,2-Trichloroethane         | ND           | U         | 100       | 100                                       | 1.0                          |
| Toluene                       | 4200         |           | 100       | 100                                       | 1.0                          |
| 2-Hexanone                    | ND           | U         | 100       | 100                                       | 1.0                          |
| Dibromochloromethane          | ND           | U         | 100       | 100                                       | 1.0                          |
| 1,2-Dibromoethane             | ND           | U         | 100       | 100                                       | 1.0                          |
| Tetrachloroethylene           | ND           | U         | 100       | 100                                       | 1.0                          |
| Chlorobenzene                 | ND           | U         | 100       | 100                                       | 1.0                          |
| Ethylbenzene                  | 1790         |           | 100       | 100                                       | 1.0                          |
| m- & p-Xylenes                | 2500         |           | 100       | 200                                       | 2.0                          |
| Bromoform                     | ND           | U         | 100       | 300                                       | 3.0                          |
| Styrene                       | ND           | U         | 100       | 100                                       | 1.0                          |
| 1,1,2,2-Tetrachloroethane     | ND           | U         | 100       | 100                                       | 1.0                          |
| o-Xylene                      | 705          |           | 100       | 100                                       | 1.0                          |
| 4-Ethyltoluene                | ND           | U         | 100       | 100                                       | 1.0                          |
| 1,3,5-Trimethylbenzene        | ND           | U         | 100       | 100                                       | 1.0                          |
| 1,2,4-Trimethylbenzene        | ND           | U         | 100       | 100                                       | 1.0                          |
| Benzyl Chloride               | ND           | U         | 100       | 500                                       | 5.0                          |
| 1,3-Dichlorobenzene           | ND           | U         | 100       | 100                                       | 1.0                          |
| 1,4-Dichlorobenzene           | ND           | U         | 100       | 100                                       | 1.0                          |
| 1,2-Dichlorobenzene           | ND           | U         | 100       | 100                                       | 1.0                          |
| 1,2,4-Trichlorobenzene        | ND           | U         | 100       | 100                                       | 1.0                          |
| Hexachlorobutadiene           | ND           | U         | 100       | 100                                       | 1.0                          |
| BFB-Surrogate Std. % Recovery | 105%         |           |           | 70-130%                                   |                              |

J- Analyte was detected. However the analyte concentration is an estimated value, which is between the Method Detection Limit (MDL) and the Reporting Limit (RL).

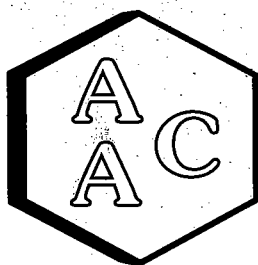
E - Estimated value, result outside linear range of instrument.

U - Compound was analyzed for, but was not detected.

!! - Estimated

  
Sucha S. Parmar, PhD  
Technical Director





# Atmospheric Analysis & Consulting, Inc.

## Laboratory Analysis Report

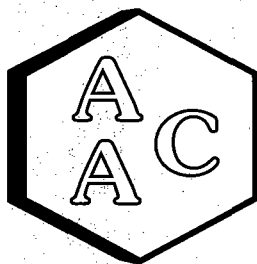
CLIENT : Shaw Environmental  
PROJECT NO : 060263  
MATRIX : AIR  
UNITS : PPB (v/v)

DATE RECEIVED : 06/09/06  
DATE REPORTED : 06/20/06

### VOLATILE ORGANIC COMPOUNDS BY EPA TO-15

| Client ID                              | A51, R2      |           |           | Sample<br>Reporting<br>Limit<br>(RLxDF's) | Method<br>Reporting<br>Limit |
|----------------------------------------|--------------|-----------|-----------|-------------------------------------------|------------------------------|
| AAC ID                                 | 060263-15747 |           |           |                                           |                              |
| Date Sampled                           | 6/7/2006     |           |           |                                           |                              |
| Date Analyzed                          | 6/17/2006    |           |           |                                           |                              |
| Can Dilution Factor                    | 1.00         |           |           |                                           |                              |
|                                        | Result       | Qualifier | Dil. Fac. |                                           |                              |
| Chlorodifluoromethane                  | 574          |           | 100       | 100                                       | 1.0                          |
| Propylene                              | 7960         |           | 1000      | 1000                                      | 1.0                          |
| Dichlorodifluoromethane                | 649          |           | 100       | 100                                       | 1.0                          |
| Chloromethane                          | ND           | U         | 100       | 100                                       | 1.0                          |
| 1,2-Dichloro-1,1,2,2-Tetrafluoroethane | ND           | U         | 100       | 100                                       | 1.0                          |
| Vinyl Chloride                         | 430          |           | 100       | 100                                       | 1.0                          |
| Methanol                               | ND           | U         | 100       | 500                                       | 5.0                          |
| 1,3-Butadiene                          | ND           | U         | 100       | 100                                       | 1.0                          |
| Bromomethane                           | ND           | U         | 100       | 100                                       | 1.0                          |
| Chloroethane                           | ND           | U         | 100       | 100                                       | 1.0                          |
| Dichlorofluoromethane                  | ND           | U         | 100       | 100                                       | 1.0                          |
| Ethanol                                | 296          |           | 100       | 200                                       | 2.0                          |
| Vinyl Bromide                          | ND           | U         | 100       | 100                                       | 1.0                          |
| Acetone                                | 983          |           | 100       | 200                                       | 2.0                          |
| Trichlorofluoromethane                 | ND           | U         | 100       | 100                                       | 1.0                          |
| Isopropyl Alcohol                      | 288          |           | 100       | 200                                       | 2.0                          |
| Acrylonitrile                          | ND           | U         | 100       | 100                                       | 1.0                          |
| 1,1-Dichloroethylene                   | ND           | U         | 100       | 100                                       | 1.0                          |
| Methylene Chloride                     | ND           | U         | 100       | 100                                       | 1.0                          |
| Allyl Chloride (Chloroprene)           | ND           | U         | 100       | 100                                       | 1.0                          |
| Carbon Disulfide                       | ND           | U         | 100       | 100                                       | 1.0                          |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane  | ND           | U         | 100       | 100                                       | 1.0                          |
| t-1,2-Dichloroethylene                 | ND           | U         | 100       | 100                                       | 1.0                          |
| 1,1-Dichloroethane                     | ND           | U         | 100       | 100                                       | 1.0                          |
| MTBE                                   | ND           | U         | 100       | 100                                       | 1.0                          |
| Vinyl Acetate                          | ND           | U         | 100       | 100                                       | 1.0                          |
| 2-Butanone (MEK)                       | 1170         |           | 100       | 100                                       | 1.0                          |
| cis-1,2- Dichloroethene                | 170          |           | 100       | 100                                       | 1.0                          |
| Hexane                                 | 454          |           | 100       | 100                                       | 1.0                          |
| Chloroform                             | ND           | U         | 100       | 100                                       | 1.0                          |
| Ethyl Acetate                          | ND           | U         | 100       | 100                                       | 1.0                          |
| Tetrahydrofuran                        | 1400         |           | 100       | 100                                       | 1.0                          |
| 1,2-Dichloroethane                     | ND           | U         | 100       | 100                                       | 1.0                          |
| 1,1,1-Trichloroethane                  | ND           | U         | 100       | 100                                       | 1.0                          |





# Atmospheric Analysis & Consulting, Inc.

## Laboratory Analysis Report

CLIENT : Shaw Environmental  
PROJECT NO : 060263  
MATRIX : AIR  
UNITS : PPB (v/v)

DATE RECEIVED : 06/09/06  
DATE REPORTED : 06/20/06

### VOLATILE ORGANIC COMPOUNDS BY EPA TO-15

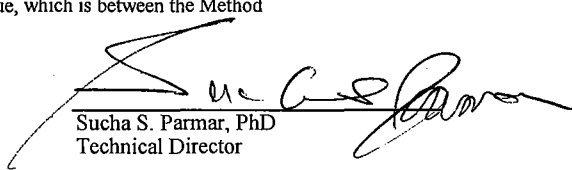
| Client ID                     | A51, R2      |           |           | Sample Reporting Limit (RLxDf's) | Method Reporting Limit |
|-------------------------------|--------------|-----------|-----------|----------------------------------|------------------------|
| AAC ID                        | 060263-15747 |           |           |                                  |                        |
| Date Sampled                  | 6/7/2006     |           |           |                                  |                        |
| Date Analyzed                 | 6/17/2006    |           |           |                                  |                        |
| Can Dilution Factor           | 1.00         |           |           |                                  |                        |
|                               | Result       | Qualifier | Dil. Fac. |                                  |                        |
| Benzene                       | 345          |           | 100       | 100                              | 1.0                    |
| Carbon Tetrachloride          | ND           | U         | 100       | 100                              | 1.0                    |
| Cyclohexane                   | 374          |           | 100       | 100                              | 1.0                    |
| 1,2-Dichloropropane           | ND           | U         | 100       | 100                              | 1.0                    |
| Bromodichloromethane          | ND           | U         | 100       | 100                              | 1.0                    |
| 1,4-Dioxane                   | ND           | U         | 100       | 100                              | 1.0                    |
| Trichloroethene               | ND           | U         | 100       | 100                              | 1.0                    |
| 2,2,4-Trimethylpentane        | 180          |           | 100       | 100                              | 1.0                    |
| Heptane                       | 858          |           | 100       | 100                              | 1.0                    |
| cis-1,3-Dichloropropene       | ND           | U         | 100       | 100                              | 1.0                    |
| 4-Methyl-2-Pentanone (MiBK)   | 162          |           | 100       | 100                              | 1.0                    |
| t-1,3-Dichloropropene         | ND           | U         | 100       | 100                              | 1.0                    |
| 1,1,2-Trichloroethane         | ND           | U         | 100       | 100                              | 1.0                    |
| Toluene                       | 4320         |           | 100       | 100                              | 1.0                    |
| 2-Hexanone                    | ND           | U         | 100       | 100                              | 1.0                    |
| Dibromochloromethane          | ND           | U         | 100       | 100                              | 1.0                    |
| 1,2-Dibromoethane             | ND           | U         | 100       | 100                              | 1.0                    |
| Tetrachloroethylene           | ND           | U         | 100       | 100                              | 1.0                    |
| Chlorobenzene                 | ND           | U         | 100       | 100                              | 1.0                    |
| Ethylbenzene                  | 1820         |           | 100       | 100                              | 1.0                    |
| m- & p-Xylenes                | 2570         |           | 100       | 200                              | 2.0                    |
| Bromoform                     | ND           | U         | 100       | 300                              | 3.0                    |
| Styrene                       | ND           | U         | 100       | 100                              | 1.0                    |
| 1,1,2,2-Tetrachloroethane     | ND           | U         | 100       | 100                              | 1.0                    |
| o-Xylene                      | 733          |           | 100       | 100                              | 1.0                    |
| 4-Ethyltoluene                | ND           | U         | 100       | 100                              | 1.0                    |
| 1,3,5-Trimethylbenzene        | ND           | U         | 100       | 100                              | 1.0                    |
| 1,2,4-Trimethylbenzene        | ND           | U         | 100       | 100                              | 1.0                    |
| Benzyl Chloride               | ND           | U         | 100       | 500                              | 5.0                    |
| 1,3-Dichlorobenzene           | ND           | U         | 100       | 100                              | 1.0                    |
| 1,4-Dichlorobenzene           | ND           | U         | 100       | 100                              | 1.0                    |
| 1,2-Dichlorobenzene           | ND           | U         | 100       | 100                              | 1.0                    |
| 1,2,4-Trichlorobenzene        | ND           | U         | 100       | 100                              | 1.0                    |
| Hexachlorobutadiene           | ND           | U         | 100       | 100                              | 1.0                    |
| BFB-Surrogate Std. % Recovery | 107%         |           |           |                                  | 70-130%                |

J- Analyte was detected. However the analyte concentration is an estimated value, which is between the Method Detection Limit (MDL) and the Reporting Limit (RL).

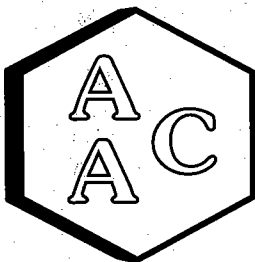
E - Estimated value, result outside linear range of instrument.

U - Compound was analyzed for, but was not detected.

!! - Estimated

  
Sucha S. Parmar, PhD  
Technical Director





# Atmospheric Analysis & Consulting, Inc.

## Laboratory Analysis Report

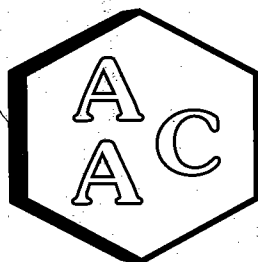
CLIENT : Shaw Environmental  
PROJECT NO : 060263  
MATRIX : AIR  
UNITS : PPB (v/v)

DATE RECEIVED : 06/09/06  
DATE REPORTED : 06/20/06

### VOLATILE ORGANIC COMPOUNDS BY EPA TO-15

| Client ID                              | A51, R2      |           |           | Sample<br>Reporting<br>Limit<br>(RLxDf's) | Method<br>Reporting<br>Limit |
|----------------------------------------|--------------|-----------|-----------|-------------------------------------------|------------------------------|
| AAC ID                                 | 060263-15747 |           |           |                                           |                              |
| Date Sampled                           | 6/7/2006     |           |           |                                           |                              |
| Date Analyzed                          | 6/17/2006    |           |           |                                           |                              |
| Can Dilution Factor                    | 1.00         |           |           |                                           |                              |
|                                        | Result       | Qualifier | Dil. Fac. |                                           |                              |
| Chlorodifluoromethane                  | 574          |           | 100       | 100                                       | 1.0                          |
| Propylene                              | 7960         |           | 1000      | 1000                                      | 1.0                          |
| Dichlorodifluoromethane                | 646          |           | 100       | 100                                       | 1.0                          |
| Chloromethane                          | ND           | U         | 100       | 100                                       | 1.0                          |
| 1,2-Dichloro-1,1,2,2-Tetrafluoroethane | ND           | U         | 100       | 100                                       | 1.0                          |
| Vinyl Chloride                         | 430          |           | 100       | 100                                       | 1.0                          |
| Methanol                               | ND           | U         | 100       | 500                                       | 5.0                          |
| 1,3-Butadiene                          | ND           | U         | 100       | 100                                       | 1.0                          |
| Bromomethane                           | ND           | U         | 100       | 100                                       | 1.0                          |
| Chloroethane                           | ND           | U         | 100       | 100                                       | 1.0                          |
| Dichlorofluoromethane                  | ND           | U         | 100       | 100                                       | 1.0                          |
| Ethanol                                | 295          |           | 100       | 200                                       | 2.0                          |
| Vinyl Bromide                          | ND           | U         | 100       | 100                                       | 1.0                          |
| Acetone                                | 1030         |           | 100       | 200                                       | 2.0                          |
| Trichlorofluoromethane                 | ND           | U         | 100       | 100                                       | 1.0                          |
| Isopropyl Alcohol                      | 279          |           | 100       | 200                                       | 2.0                          |
| Acrylonitrile                          | ND           | U         | 100       | 100                                       | 1.0                          |
| 1,1-Dichloroethylene                   | ND           | U         | 100       | 100                                       | 1.0                          |
| Methylene Chloride                     | ND           | U         | 100       | 100                                       | 1.0                          |
| Allyl Chloride (Chloroprene)           | ND           | U         | 100       | 100                                       | 1.0                          |
| Carbon Disulfide                       | ND           | U         | 100       | 100                                       | 1.0                          |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane  | ND           | U         | 100       | 100                                       | 1.0                          |
| t-1,2-Dichloroethylene                 | ND           | U         | 100       | 100                                       | 1.0                          |
| 1,1-Dichloroethane                     | ND           | U         | 100       | 100                                       | 1.0                          |
| MTBE                                   | ND           | U         | 100       | 100                                       | 1.0                          |
| Vinyl Acetate                          | ND           | U         | 100       | 100                                       | 1.0                          |
| 2-Butanone (MEK)                       | 1190         |           | 100       | 100                                       | 1.0                          |
| cis-1,2- Dichloroethene                | 172          |           | 100       | 100                                       | 1.0                          |
| Hexane                                 | 463          |           | 100       | 100                                       | 1.0                          |
| Chloroform                             | ND           | U         | 100       | 100                                       | 1.0                          |
| Ethyl Acetate                          | ND           | U         | 100       | 100                                       | 1.0                          |
| Tetrahydrofuran                        | 1450         |           | 100       | 100                                       | 1.0                          |
| 1,2-Dichloroethane                     | ND           | U         | 100       | 100                                       | 1.0                          |
| 1,1,1-Trichloroethane                  | ND           | U         | 100       | 100                                       | 1.0                          |





# Atmospheric Analysis & Consulting, Inc.

## Laboratory Analysis Report

CLIENT : Shaw Environmental  
PROJECT NO : 060263  
MATRIX : AIR  
UNITS : PPB (v/v)

DATE RECEIVED : 06/09/06  
DATE REPORTED : 06/20/06

### VOLATILE ORGANIC COMPOUNDS BY EPA TO-15

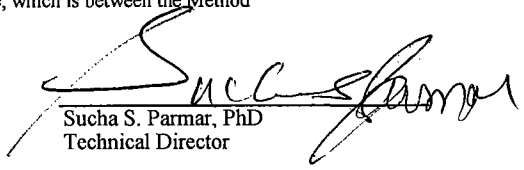
| Client ID                     | A51, R2      |           |           | Sample<br>Reporting<br>Limit<br>(RLxDf's) | Method<br>Reporting<br>Limit |
|-------------------------------|--------------|-----------|-----------|-------------------------------------------|------------------------------|
| AAC ID                        | 060263-15747 |           |           |                                           |                              |
| Date Sampled                  | 6/7/2006     |           |           |                                           |                              |
| Date Analyzed                 | 6/17/2006    |           |           |                                           |                              |
| Can Dilution Factor           | 1.00         |           |           |                                           |                              |
|                               | Result       | Qualifier | Dil. Fac. |                                           |                              |
| Benzene                       | 336          |           | 100       | 100                                       | 1.0                          |
| Carbon Tetrachloride          | ND           | U         | 100       | 100                                       | 1.0                          |
| Cyclohexane                   | 362          |           | 100       | 100                                       | 1.0                          |
| 1,2-Dichloropropane           | ND           | U         | 100       | 100                                       | 1.0                          |
| Bromodichloromethane          | ND           | U         | 100       | 100                                       | 1.0                          |
| 1,4-Dioxane                   | ND           | U         | 100       | 100                                       | 1.0                          |
| Trichloroethene               | ND           | U         | 100       | 100                                       | 1.0                          |
| 2,2,4-Trimethylpentane        | 174          |           | 100       | 100                                       | 1.0                          |
| Heptane                       | 844          |           | 100       | 100                                       | 1.0                          |
| cis-1,3-Dichloropropene       | ND           | U         | 100       | 100                                       | 1.0                          |
| 4-Methyl-2-Pentanone (MiBK)   | 157          |           | 100       | 100                                       | 1.0                          |
| t-1,3-Dichloropropene         | ND           | U         | 100       | 100                                       | 1.0                          |
| 1,1,2-Trichloroethane         | ND           | U         | 100       | 100                                       | 1.0                          |
| Toluene                       | 4260         |           | 100       | 100                                       | 1.0                          |
| 2-Hexanone                    | ND           | U         | 100       | 100                                       | 1.0                          |
| Dibromochloromethane          | ND           | U         | 100       | 100                                       | 1.0                          |
| 1,2-Dibromoethane             | ND           | U         | 100       | 100                                       | 1.0                          |
| Tetrachloroethylene           | ND           | U         | 100       | 100                                       | 1.0                          |
| Chlorobenzene                 | ND           | U         | 100       | 100                                       | 1.0                          |
| Ethylbenzene                  | 1800         |           | 100       | 100                                       | 1.0                          |
| m- & p-Xylenes                | 2540         |           | 100       | 200                                       | 2.0                          |
| Bromoform                     | ND           | U         | 100       | 300                                       | 3.0                          |
| Styrene                       | ND           | U         | 100       | 100                                       | 1.0                          |
| 1,1,2,2-Tetrachloroethane     | ND           | U         | 100       | 100                                       | 1.0                          |
| o-Xylene                      | 719          |           | 100       | 100                                       | 1.0                          |
| 4-Ethyltoluene                | ND           | U         | 100       | 100                                       | 1.0                          |
| 1,3,5-Trimethylbenzene        | ND           | U         | 100       | 100                                       | 1.0                          |
| 1,2,4-Trimethylbenzene        | ND           | U         | 100       | 100                                       | 1.0                          |
| Benzyl Chloride               | ND           | U         | 100       | 500                                       | 5.0                          |
| 1,3-Dichlorobenzene           | ND           | U         | 100       | 100                                       | 1.0                          |
| 1,4-Dichlorobenzene           | ND           | U         | 100       | 100                                       | 1.0                          |
| 1,2-Dichlorobenzene           | ND           | U         | 100       | 100                                       | 1.0                          |
| 1,2,4-Trichlorobenzene        | ND           | U         | 100       | 100                                       | 1.0                          |
| Hexachlorobutadiene           | ND           | U         | 100       | 100                                       | 1.0                          |
| BFB-Surrogate Std. % Recovery | 106%         |           |           | 70-130%                                   |                              |

J- Analyte was detected. However the analyte concentration is an estimated value, which is between the Method Detection Limit (MDL) and the Reporting Limit (RL).

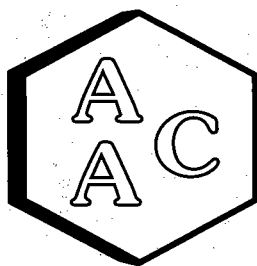
E - Estimated value, result outside linear range of instrument.

U - Compound was analyzed for, but was not detected.

!! - Estimated

  
Sucha S. Parmar, PhD  
Technical Director





# Atmospheric Analysis & Consulting, Inc.

## Method Blank Analysis Report

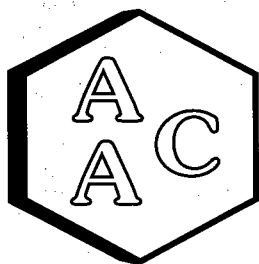
MATRIX : AIR  
UNITS : ppbv

ANALYSIS DATE : 06/12/06  
REPORT DATE : 06/16/06

### VOLATILE ORGANIC COMPOUNDS BY EPA TO-14/TO-15

| Client ID<br>AAC ID                    | Method Blank<br>MB 061206 | RL  |
|----------------------------------------|---------------------------|-----|
| Chlorodifluoromethane*                 | <RL                       | 1.0 |
| Propylene*                             | <RL                       | 1.0 |
| DiCIDIEMethane*                        | <RL                       | 1.0 |
| CHLOROMETHANE*                         | <RL                       | 1.0 |
| 1,2 DiCl-1,1,2,2-TetraFEthane*         | <RL                       | 1.0 |
| VINYL CHLORIDE*                        | <RL                       | 1.0 |
| Methanol*                              | <RL                       | 5.0 |
| 1,3-Butadiene*                         | <RL                       | 1.0 |
| BROMOMETHANE*                          | <RL                       | 1.0 |
| CHLOROETHANE*                          | <RL                       | 1.0 |
| Dichlorofluoromethane                  | <RL                       | 1.0 |
| Ethanol*                               | <RL                       | 2.0 |
| Vinyl Bromide*                         | <RL                       | 1.0 |
| Acetone*                               | <RL                       | 2.0 |
| TRICHLOROFLUOROMETHANE*                | <RL                       | 1.0 |
| Isopropyl Alcohol*                     | <RL                       | 2.0 |
| Acrylonitrile*                         | <RL                       | 1.0 |
| 1,1 DICHLOROETHENE*                    | <RL                       | 1.0 |
| METHYLENE CHLORIDE*                    | <RL                       | 1.0 |
| Allyl CHLORIDE*                        | <RL                       | 1.0 |
| Carbon disulfide*                      | <RL                       | 1.0 |
| 1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE* | <RL                       | 1.0 |
| trans-1,2- DICHLOROETHYLENE*           | <RL                       | 1.0 |
| 1,1- DICHLOROETHANE*                   | <RL                       | 1.0 |
| MTBE*                                  | <RL                       | 1.0 |
| Vinyl Acetate*                         | <RL                       | 1.0 |
| MEK*                                   | <RL                       | 1.0 |
| cis-1,2- DICHLOROETHYLENE*             | <RL                       | 1.0 |
| Hexane*                                | <RL                       | 1.0 |
| CHLOROFORM*                            | <RL                       | 1.0 |
| Ethyl Acetate*                         | <RL                       | 1.0 |
| Tetrahydrofuran*                       | <RL                       | 1.0 |
| 1,2-DICHLOROETHANE*                    | <RL                       | 1.0 |
| 1,1,1-TRICHLOROETHANE*                 | <RL                       | 1.0 |
| BENZENE**                              | <RL                       | 1.0 |
| CARBON TETRACHLORIDE**                 | <RL                       | 1.0 |
| Cyclohexane**                          | <RL                       | 1.0 |
| 1,2-DICHLOROPROPANE**                  | <RL                       | 1.0 |
| Bromodichloromethane**                 | <RL                       | 1.0 |
| 1,4-Dioxane**                          | <RL                       | 1.0 |
| TRICHLOROETHENE**                      | <RL                       | 1.0 |
| 2,2,4-Trimethylpentane**               | <RL                       | 1.0 |
| Heptane**                              | <RL                       | 1.0 |





## Atmospheric Analysis & Consulting, Inc.

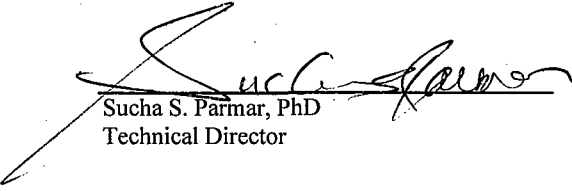
### Method Blank Analysis Report

MATRIX : AIR ANALYSIS DATE : 06/12/06  
UNITS : ppbv REPORT DATE : 06/16/06

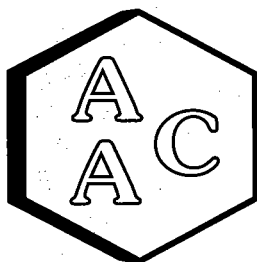
#### VOLATILE ORGANIC COMPOUNDS BY EPA TO-14/TO-15

| Client ID<br>AAC ID                | Method Blank<br>MB 061206 | RL  |
|------------------------------------|---------------------------|-----|
| cis- 1,3 DICHLOROPROPENE**         | <RL                       | 1.0 |
| MIBK**                             | <RL                       | 1.0 |
| trans 1,3 DICHLOROPROPENE**        | <RL                       | 1.0 |
| 1,1,2-TRICHLOROETHANE**            | <RL                       | 1.0 |
| TOLUENE**                          | <RL                       | 1.0 |
| 2-Hexanone**                       | <RL                       | 1.0 |
| Dibromochloromethane**             | <RL                       | 1.0 |
| 1,2 DIBROMOETHANE**                | <RL                       | 1.0 |
| TETRACHLOROETHYLENE**              | <RL                       | 1.0 |
| CHLOROBENZENE***                   | <RL                       | 1.0 |
| ETHYLBENZENE***                    | <RL                       | 1.0 |
| m-, & p- XYLENES***                | <RL                       | 2.0 |
| Bromoform***                       | <RL                       | 3.0 |
| STYRENE***                         | <RL                       | 1.0 |
| 1,1, 2,2- TETRACHLOROETHANE***     | <RL                       | 1.0 |
| o- XYLENE***                       | <RL                       | 1.0 |
| Ethyltoluene***                    | <RL                       | 1.0 |
| 1,3,5- TRIMETHYLBENZENE***         | <RL                       | 1.0 |
| 1,2,4- TRIMETHYLBENZENE***         | <RL                       | 1.0 |
| Benzyl Chloride***                 | <RL                       | 5.0 |
| 1,3- DICHLOROBENZENE***            | <RL                       | 1.0 |
| 1,4- DICHLOROBENZENE***            | <RL                       | 1.0 |
| 1,2-DICHLOROBENZENE***             | <RL                       | 1.0 |
| 1,2,4 TRICHLOROBENZENE***          | <RL                       | 1.0 |
| HEXACHLOROBUTADIENE***             | <RL                       | 1.0 |
| <b>System Monitoring Compounds</b> |                           |     |
| BFB-Surrogate Std. % Recovery      | 97%                       | --  |

RL - Reporting Limit

  
Sucha S. Parmar, PhD  
Technical Director





# Atmospheric Analysis & Consulting, Inc.

## Quality Control/Quality Assurance Report

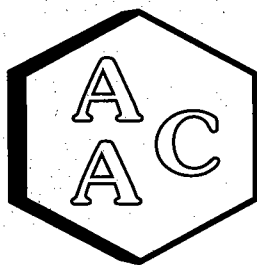
AAC ID : 060264-15752  
MATRIX : Air

DATE ANALYZED : 06/12/06  
DATE REPORTED : 06/16/06  
UNITS : ppbv

### TO-14/TO-15 Duplicate Analysis

| Compound                               | Sample Conc | Duplicate Conc | % RPD |
|----------------------------------------|-------------|----------------|-------|
| Chlorodifluoromethane*                 | <RL         | <RL            | 0.0   |
| Propylene*                             | <RL         | <RL            | 0.0   |
| DiCIDIEMethane*                        | <RL         | <RL            | 0.0   |
| CHLOROMETHANE*                         | <RL         | <RL            | 0.0   |
| 1,2 DiCl-1,1,2,2-TetraF Ethane*        | <RL         | <RL            | 0.0   |
| VINYL CHLORIDE*                        | <RL         | <RL            | 0.0   |
| Methanol*                              | <RL         | <RL            | 0.0   |
| 1,3-Butadiene*                         | <RL         | <RL            | 0.0   |
| BROMOMETHANE*                          | <RL         | <RL            | 0.0   |
| CHLOROETHANE*                          | <RL         | <RL            | 0.0   |
| Dichlorofluoromethane                  | <RL         | <RL            | 0.0   |
| Ethanol*                               | <RL         | <RL            | 0.0   |
| Vinyl Bromide*                         | <RL         | <RL            | 0.0   |
| Acetone*                               | 5.79        | 5.50           | 5.1   |
| TRICHLOROFLUOROMETHANE*                | <RL         | <RL            | 0.0   |
| Isopropyl Alcohol*                     | <RL         | <RL            | 0.0   |
| Acrylonitrile*                         | <RL         | <RL            | 0.0   |
| 1,1 DICHLOROETHENE*                    | <RL         | <RL            | 0.0   |
| METHYLENE CHLORIDE*                    | <RL         | <RL            | 0.0   |
| Allyl CHLORIDE*                        | <RL         | <RL            | 0.0   |
| Carbon disulfide*                      | <RL         | <RL            | 0.0   |
| 1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE* | <RL         | <RL            | 0.0   |
| trans-1,2- DICHLOROETHYLENE*           | <RL         | <RL            | 0.0   |
| 1,1- DICHLOROETHANE*                   | <RL         | <RL            | 0.0   |
| MTBE*                                  | <RL         | <RL            | 0.0   |
| Vinyl Acetate*                         | <RL         | <RL            | 0.0   |
| MEK*                                   | <RL         | <RL            | 0.0   |
| cis-1,2- DICHLOROETHYLENE*             | <RL         | <RL            | 0.0   |
| Hexane*                                | <RL         | <RL            | 0.0   |
| CHLOROFORM*                            | <RL         | <RL            | 0.0   |
| Ethyl Acetate*                         | <RL         | <RL            | 0.0   |
| Tetrahydrofuran*                       | <RL         | <RL            | 0.0   |
| 1,2-DICHLOROETHANE*                    | <RL         | <RL            | 0.0   |
| 1,1,1-TRICHLOROETHANE*                 | <RL         | <RL            | 0.0   |
| BENZENE**                              | <RL         | <RL            | 0.0   |
| CARBON TETRACHLORIDE**                 | <RL         | <RL            | 0.0   |





# Atmospheric Analysis & Consulting, Inc.

## Quality Control/Quality Assurance Report

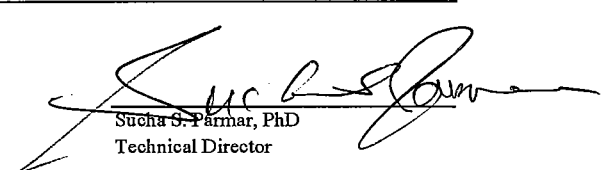
AAC ID : 060264-15752  
MATRIX : Air

DATE ANALYZED : 06/12/06  
DATE REPORTED : 06/16/06  
UNITS : ppbv

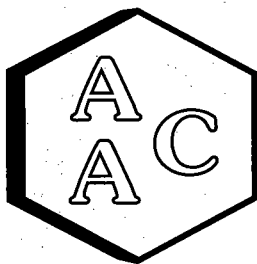
### TO-14/TO-15 Duplicate Analysis

| Compound                      | Sample Conc | Duplicate Conc | % RPD |
|-------------------------------|-------------|----------------|-------|
| Cyclohexane**                 | <RL         | <RL            | 0.0   |
| 1,2-DICHLOROPROPANE**         | <RL         | <RL            | 0.0   |
| Bromodichloromethane**        | <RL         | <RL            | 0.0   |
| 1,4-Dioxane**                 | <RL         | <RL            | 0.0   |
| TRICHLOROETHENE**             | <RL         | <RL            | 0.0   |
| 2,2,4-Trimethylpentane**      | <RL         | <RL            | 0.0   |
| Heptane**                     | <RL         | <RL            | 0.0   |
| cis- 1,3 DICHLOROPROPENE**    | <RL         | <RL            | 0.0   |
| MIBK**                        | <RL         | <RL            | 0.0   |
| trans 1,3 DICHLOROPROPENE**   | <RL         | <RL            | 0.0   |
| 1,1,2-TRICHLOROETHANE**       | <RL         | <RL            | 0.0   |
| TOLUENE**                     | <RL         | <RL            | 0.0   |
| 2-Hexanone**                  | <RL         | <RL            | 0.0   |
| Dibromochloromethane**        | <RL         | <RL            | 0.0   |
| 1,2 DIBROMOETHANE**           | <RL         | <RL            | 0.0   |
| TETRACHLOROETHYLENE**         | <RL         | <RL            | 0.0   |
| CHLOROBENZENE***              | <RL         | <RL            | 0.0   |
| ETHYLBENZENE***               | <RL         | <RL            | 0.0   |
| m-, & p- XYLENES***           | <RL         | <RL            | 0.0   |
| Bromoform***                  | <RL         | <RL            | 0.0   |
| STYRENE***                    | <RL         | <RL            | 0.0   |
| 1,1,2,2-TETRACHLORETHANE***   | <RL         | <RL            | 0.0   |
| o- XYLENE***                  | <RL         | <RL            | 0.0   |
| Ethyltoluene***               | <RL         | <RL            | 0.0   |
| 1,3,5- TRIMETHYLBENZENE***    | <RL         | <RL            | 0.0   |
| 1,2,4- TRIMETHYLBENZENE***    | <RL         | <RL            | 0.0   |
| Benzyl Chloride***            | <RL         | <RL            | 0.0   |
| 1,3- DICHLOROBENZENE***       | <RL         | <RL            | 0.0   |
| 1,4- DICHLOROBENZENE***       | <RL         | <RL            | 0.0   |
| 1,2-DICHLOROBENZENE***        | <RL         | <RL            | 0.0   |
| 1,2,4 TRICHLOROBENZENE***     | <RL         | <RL            | 0.0   |
| Hexachlorobutadiene***        | <RL         | <RL            | 0.0   |
| System Monitoring Compounds   |             |                |       |
| BFB-Surrogate Std. % Recovery | 95          | 98             | 2.8   |

RL - Reporting Limit

  
Sucha G. Parmar, PhD  
Technical Director





## Atmospheric Analysis & Consulting, Inc.

### Quality Control/Quality Assurance Report

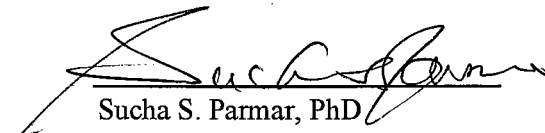
CLIENT ID : Laboratory Control Spike      DATE ANALYZED : 06/17/06  
AAC ID : LCS/LCSD      DATE REPORTED : 06/19/06  
MEDIA : Air      UNITS : ppbv

### TO-14/15 Laboratory Control Spike Recovery

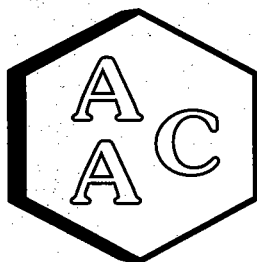
| Compound             | Sample Conc. | Spike Added | Spike Res | Dup Spike Res | Spike % Rec * | Spike Dup % Rec * | RPD** % |
|----------------------|--------------|-------------|-----------|---------------|---------------|-------------------|---------|
| 1,1-DICHLOROETHYLENE | 0.0          | 10.00       | 8.28      | 8.61          | 83            | 86                | 3.9     |
| METHYLENECHLORIDE    | 0.0          | 10.00       | 10.93     | 11.19         | 109           | 112               | 2.4     |
| BENZENE              | 0.0          | 10.00       | 10.61     | 11.02         | 106           | 110               | 3.8     |
| TRICHLOROETHENE      | 0.0          | 10.00       | 11.04     | 11.51         | 110           | 115               | 4.2     |
| TOLUENE              | 0.0          | 10.00       | 10.50     | 10.99         | 105           | 110               | 4.6     |
| TETRACHLOROETHYLENE  | 0.0          | 10.00       | 10.27     | 10.88         | 103           | 109               | 5.8     |
| CHLOROBENZENE        | 0.0          | 10.00       | 10.60     | 11.09         | 106           | 111               | 4.5     |
| ETHYLBENZENE         | 0.0          | 10.00       | 10.64     | 11.20         | 106           | 112               | 5.1     |
| m-, & p- XYLENES     | 0.0          | 20.00       | 21.57     | 23.25         | 108           | 116               | 7.5     |
| o- XYLENE            | 0.0          | 10.00       | 11.00     | 11.66         | 110           | 117               | 5.8     |

\* Must be 70-130%

\*\* Must be < 25%

  
Sucha S. Parmar, PhD  
Technical Director





## Atmospheric Analysis & Consulting, Inc.

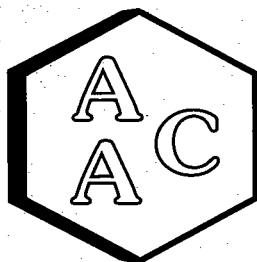
### Method Blank Analysis Report

MATRIX : AIR ANALYSIS DATE : 06/17/06  
UNITS : ppbv REPORT DATE : 06/19/06

#### VOLATILE ORGANIC COMPOUNDS BY EPA TO-14/TO-15

| Client ID                              | Method Blank | RL  |
|----------------------------------------|--------------|-----|
| AAC ID                                 | MB 061706    |     |
| Chlorodifluoromethane*                 | <RL          | 1.0 |
| Propylene*                             | <RL          | 1.0 |
| DiCIDIFMethane*                        | <RL          | 1.0 |
| CHLOROMETHANE*                         | <RL          | 1.0 |
| 1,2 DiCl-1,1,2,2-TetraFEthane*         | <RL          | 1.0 |
| VINYL CHLORIDE*                        | <RL          | 1.0 |
| Methanol*                              | <RL          | 5.0 |
| 1,3-Butadiene*                         | <RL          | 1.0 |
| BROMOMETHANE*                          | <RL          | 1.0 |
| CHLOROETHANE*                          | <RL          | 1.0 |
| Dichlorofluoromethane                  | <RL          | 1.0 |
| Ethanol*                               | <RL          | 2.0 |
| Vinyl Bromide*                         | <RL          | 1.0 |
| Acetone*                               | <RL          | 2.0 |
| TRICHLOROFLUOROMETHANE*                | <RL          | 1.0 |
| Isopropyl Alcohol*                     | <RL          | 2.0 |
| Acrylonitrile*                         | <RL          | 1.0 |
| 1,1 DICHLOROETHENE*                    | <RL          | 1.0 |
| METHYLENE CHLORIDE*                    | <RL          | 1.0 |
| Allyl CHLORIDE*                        | <RL          | 1.0 |
| Carbon disulfide*                      | <RL          | 1.0 |
| 1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE* | <RL          | 1.0 |
| trans-1,2- DICHLOROETHYLENE*           | <RL          | 1.0 |
| 1,1- DICHLOROETHANE*                   | <RL          | 1.0 |
| MTBE*                                  | <RL          | 1.0 |
| Vinyl Acetate*                         | <RL          | 1.0 |
| MEK*                                   | <RL          | 1.0 |
| cis-1,2- DICHLOROETHYLENE*             | <RL          | 1.0 |
| Hexane*                                | <RL          | 1.0 |
| CHLOROFORM*                            | <RL          | 1.0 |
| Ethyl Acetate*                         | <RL          | 1.0 |
| Tetrahydrofuran*                       | <RL          | 1.0 |
| 1,2-DICHLOROETHANE*                    | <RL          | 1.0 |
| 1,1,1-TRICHLOROETHANE*                 | <RL          | 1.0 |
| BENZENE**                              | <RL          | 1.0 |
| CARBON TETRACHLORIDE**                 | <RL          | 1.0 |
| Cyclohexane**                          | <RL          | 1.0 |
| 1,2-DICHLOROPROPANE**                  | <RL          | 1.0 |
| Bromodichloromethane**                 | <RL          | 1.0 |
| 1,4-Dioxane**                          | <RL          | 1.0 |
| TRICHLOROETHENE**                      | <RL          | 1.0 |
| 2,2,4-Trimethylpentane**               | <RL          | 1.0 |
| Heptane**                              | <RL          | 1.0 |





## Atmospheric Analysis & Consulting, Inc.

### Method Blank Analysis Report

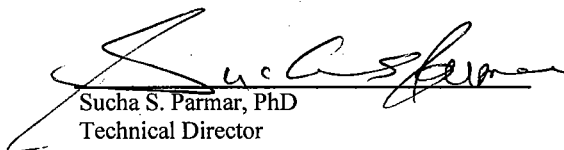
MATRIX : AIR  
UNITS : ppbv

ANALYSIS DATE : 06/17/06  
REPORT DATE : 06/19/06

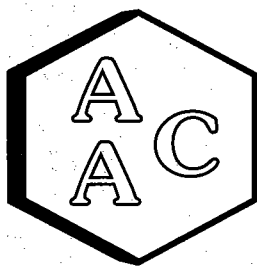
#### VOLATILE ORGANIC COMPOUNDS BY EPA TO-14/TO-15

| Client ID<br>AAC ID            | Method Blank<br>MB 061706 | RL  |
|--------------------------------|---------------------------|-----|
| cis- 1,3 DICHLOROPROPENE**     | <RL                       | 1.0 |
| MiBK**                         | <RL                       | 1.0 |
| trans 1,3 DICHLOROPROPENE**    | <RL                       | 1.0 |
| 1,1,2-TRICHLOROETHANE**        | <RL                       | 1.0 |
| TOLUENE**                      | <RL                       | 1.0 |
| 2-Hexanone**                   | <RL                       | 1.0 |
| Dibromochloromethane**         | <RL                       | 1.0 |
| 1,2 DIBROMOETHANE**            | <RL                       | 1.0 |
| TETRACHLOROETHYLENE**          | <RL                       | 1.0 |
| CHLORO BENZENE***              | <RL                       | 1.0 |
| ETHYL BENZENE***               | <RL                       | 1.0 |
| m-, & p- XYLENES***            | <RL                       | 2.0 |
| Bromoform***                   | <RL                       | 3.0 |
| STYRENE***                     | <RL                       | 1.0 |
| 1,1, 2,2- TETRACHLOROETHANE*** | <RL                       | 1.0 |
| o- XYLENE***                   | <RL                       | 1.0 |
| Ethyltoluene***                | <RL                       | 1.0 |
| 1,3,5- TRIMETHYLBENZENE***     | <RL                       | 1.0 |
| 1,2,4- TRIMETHYLBENZENE***     | <RL                       | 1.0 |
| Benzyl Chloride***             | <RL                       | 5.0 |
| 1,3- DICHLOROBENZENE***        | <RL                       | 1.0 |
| 1,4- DICHLOROBENZENE***        | <RL                       | 1.0 |
| 1,2-DICHLOROBENZENE***         | <RL                       | 1.0 |
| 1,2,4 TRICHLOROBENZENE***      | <RL                       | 1.0 |
| HEXACHLOROBUTADIENE***         | <RL                       | 1.0 |
| System Monitoring Compounds    |                           |     |
| BFB-Surrogate Std. % Recovery  | 96%                       | --  |

RL - Reporting Limit

  
Sucha S. Parmar, PhD  
Technical Director





# Atmospheric Analysis & Consulting, Inc.

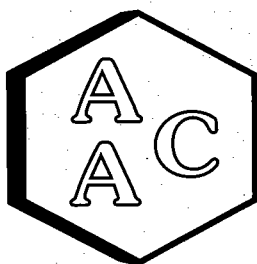
## Quality Control/Quality Assurance Report

AAC ID : 060283-15834      DATE ANALYZED : 06/17/06  
 MATRIX : Air      DATE REPORTED : 06/19/06  
 UNITS : ppbv

### TO-14/TO-15 Duplicate Analysis

| Compound                               | Sample Conc | Duplicate Conc | % RPD |
|----------------------------------------|-------------|----------------|-------|
| Chlorodifluoromethane*                 | <RL         | <RL            | 0.0   |
| Propylene*                             | <RL         | <RL            | 0.0   |
| DiClDiFMethane*                        | <RL         | <RL            | 0.0   |
| CHLOROMETHANE*                         | <RL         | <RL            | 0.0   |
| 1,2 DiCl-1,1,2,2-TetraFEthane*         | <RL         | <RL            | 0.0   |
| VINYL CHLORIDE*                        | <RL         | <RL            | 0.0   |
| Methanol*                              | 11.30       | 9.78           | 14.4  |
| 1,3-Butadiene*                         | <RL         | <RL            | 0.0   |
| BROMOMETHANE*                          | <RL         | <RL            | 0.0   |
| CHLOROETHANE*                          | <RL         | <RL            | 0.0   |
| Dichlorofluoromethane                  | <RL         | <RL            | 0.0   |
| Ethanol*                               | 4.92        | 4.18           | 16.3  |
| Vinyl Bromide*                         | <RL         | <RL            | 0.0   |
| Acetone*                               | 7.90        | 7.82           | 1.0   |
| TRICHLOROFLUOROMETHANE*                | <RL         | <RL            | 0.0   |
| Isopropyl Alcohol*                     | <RL         | <RL            | 0.0   |
| Acrylonitrile*                         | <RL         | <RL            | 0.0   |
| 1,1 DICHLOROETHENE*                    | <RL         | <RL            | 0.0   |
| METHYLENE CHLORIDE*                    | 6.43        | 6.19           | 3.8   |
| Allyl CHLORIDE*                        | <RL         | <RL            | 0.0   |
| Carbon disulfide*                      | 1.50        | 1.43           | 4.8   |
| 1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE* | <RL         | <RL            | 0.0   |
| trans-1,2- DICHLOROETHYLENE*           | <RL         | <RL            | 0.0   |
| 1,1- DICHLOROETHANE*                   | <RL         | <RL            | 0.0   |
| MTBE*                                  | <RL         | <RL            | 0.0   |
| Vinyl Acetate*                         | <RL         | <RL            | 0.0   |
| MEK*                                   | <RL         | <RL            | 0.0   |
| cis-1,2- DICHLOROETHYLENE*             | <RL         | <RL            | 0.0   |
| Hexane*                                | 2.30        | 2.32           | 0.9   |
| CHLOROFORM*                            | <RL         | <RL            | 0.0   |
| Ethyl Acetate*                         | <RL         | <RL            | 0.0   |
| Tetrahydrofuran*                       | <RL         | <RL            | 0.0   |
| 1,2-DICHLOROETHANE*                    | <RL         | <RL            | 0.0   |
| 1,1,1-TRICHLOROETHANE*                 | <RL         | <RL            | 0.0   |
| BENZENE**                              | <RL         | <RL            | 0.0   |
| CARBON TETRACHLORIDE**                 | <RL         | <RL            | 0.0   |





# Atmospheric Analysis & Consulting, Inc.

## Quality Control/Quality Assurance Report

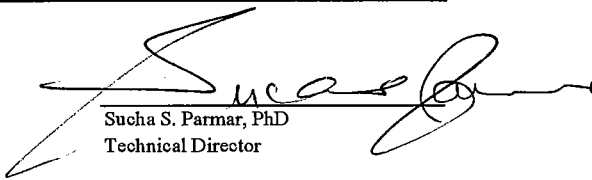
AAC ID : 060283-15834  
MATRIX : Air

DATE ANALYZED : 06/17/06  
DATE REPORTED : 06/19/06  
UNITS : ppbv

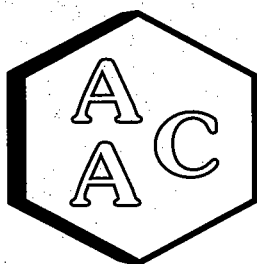
### TO-14/TO-15 Duplicate Analysis

| Compound                      | Sample Conc | Duplicate Conc | % RPD |
|-------------------------------|-------------|----------------|-------|
| Cyclohexane**                 | <RL         | <RL            | 0.0   |
| 1,2-DICHLOROPROPANE**         | <RL         | <RL            | 0.0   |
| Bromodichloromethane**        | <RL         | <RL            | 0.0   |
| 1,4-Dioxane**                 | <RL         | <RL            | 0.0   |
| TRICHLOROETHENE**             | <RL         | <RL            | 0.0   |
| 2,2,4-Trimethylpentane**      | <RL         | <RL            | 0.0   |
| Heptane**                     | <RL         | <RL            | 0.0   |
| cis- 1,3 DICHLOROPROPENE**    | <RL         | <RL            | 0.0   |
| MIBK**                        | <RL         | <RL            | 0.0   |
| trans 1,3 DICHLOROPROPENE**   | <RL         | <RL            | 0.0   |
| 1,1,2-TRICHLOROETHANE**       | <RL         | <RL            | 0.0   |
| TOLUENE**                     | 2.40        | 2.43           | 1.2   |
| 2-Hexanone**                  | <RL         | <RL            | 0.0   |
| Dibromochloromethane**        | <RL         | <RL            | 0.0   |
| 1,2 DIBROMOETHANE**           | <RL         | <RL            | 0.0   |
| TETRACHLOROETHYLENE**         | <RL         | <RL            | 0.0   |
| CHLOROBENZENE***              | <RL         | <RL            | 0.0   |
| ETHYLBENZENE***               | <RL         | <RL            | 0.0   |
| m-, & p- XYLENES***           | <RL         | <RL            | 0.0   |
| Bromoform***                  | <RL         | <RL            | 0.0   |
| STYRENE***                    | <RL         | <RL            | 0.0   |
| 1,1,2,2-TETRACHLORETHANE***   | <RL         | <RL            | 0.0   |
| o- XYLENE***                  | <RL         | <RL            | 0.0   |
| Ethyltoluene***               | <RL         | <RL            | 0.0   |
| 1,3,5- TRIMETHYLBENZENE***    | <RL         | <RL            | 0.0   |
| 1,2,4- TRIMETHYLBENZENE***    | <RL         | <RL            | 0.0   |
| Benzyl Chloride***            | <RL         | <RL            | 0.0   |
| 1,3- DICHLOROBENZENE***       | <RL         | <RL            | 0.0   |
| 1,4- DICHLOROBENZENE***       | <RL         | <RL            | 0.0   |
| 1,2-DICHLOROBENZENE***        | <RL         | <RL            | 0.0   |
| 1,2,4 TRICHLOROBENZENE***     | <RL         | <RL            | 0.0   |
| Hexachlorobutadiene***        | <RL         | <RL            | 0.0   |
| System Monitoring Compounds   |             |                |       |
| BFB-Surrogate Std. % Recovery | 99          | 98             | 0.8   |

RL - Reporting Limit

  
Sucha S. Parmar, PhD  
Technical Director





# Atmospheric Analysis & Consulting, Inc.

ANALYSIS DATE : 06/17/06

INSTRUMENT ID

: GC/MS-02

ANALYST : JIG

STD ID

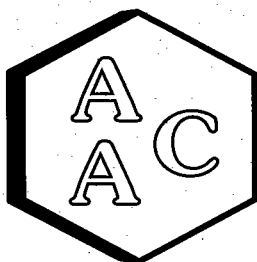
: PS061406-09

## VOLATILE ORGANIC COMPOUNDS BY EPA METHOD TO-14/TO-15

Continuing Calibration Verification of the 06/16/06 Calibration

| Compounds                      | Conc | Daily Conc | %REC |
|--------------------------------|------|------------|------|
| 4-BFB (surrogate standard)***  | 10   | 10.94      | 109  |
| Chlorodifluoromethane*         | 10   | 10.76      | 108  |
| Propylene*                     | 10   | 10.53      | 105  |
| DiCIDIFMethane*                | 10   | 11.78      | 118  |
| CHLOROMETHANE*                 | 10   | 11.05      | 111  |
| 1,2 DiCl-1,1,2,2-TetraFEthane* | 10   | 8.90       | 89   |
| VINYL CHLORIDE*                | 10   | 11.03      | 110  |
| Methanol*                      | 10   | 11.23      | 112  |
| 1,3-Butadiene*                 | 10   | 11.41      | 114  |
| BROMOMETHANE*                  | 10   | 11.03      | 110  |
| CHLOROETHANE*                  | 10   | 11.01      | 110  |
| Dichlorofluoromethane*         | 10   | 11.59      | 116  |
| Ethanol*                       | 10   | 10.91      | 109  |
| Vinyl Bromide*                 | 10   | 8.70       | 87   |
| Acetone*                       | 10   | 11.06      | 111  |
| TRICHLOROFLUOROMETHANE*        | 10   | 11.49      | 115  |
| Isopropanol*                   | 10   | 10.81      | 108  |
| Acrylonitrile*                 | 10   | 10.68      | 107  |
| 1,1 DICHLOROETHENE*            | 10   | 8.28       | 83   |
| METHYLENE CHLORIDE*            | 10   | 10.93      | 109  |
| Allyl CHLORIDE*                | 10   | 8.10       | 81   |
| Carbon disulfide*              | 10   | 9.04       | 90   |
| 1,1,2-TRICHLORO-1,2,2-TRIFLUO  | 10   | 9.43       | 94   |
| trans-1,2- DICHLOROETHYLENE*   | 10   | 8.15       | 82   |
| 1,1- DICHLOROETHANE*           | 10   | 7.85       | 79   |
| MTBE*                          | 10   | 7.70       | 77   |
| Vinyl Acetate*                 | 10   | 7.72       | 77   |
| MEK*                           | 10   | 9.69       | 97   |
| cis-1,2- DICHLOROETHYLENE*     | 10   | 10.37      | 104  |
| Hexane*                        | 10   | 10.23      | 102  |
| CHLOROFORM*                    | 10   | 10.14      | 101  |
| Ethyl Acetate*                 | 10   | 10.25      | 103  |
| Tetrahydrofuran*               | 10   | 9.30       | 93   |
| 1,2-DICHLOROETHANE*            | 10   | 9.48       | 95   |
| 1,1,1-TRICHLOROETHANE*         | 10   | 8.91       | 89   |





# Atmospheric Analysis & Consulting, Inc.

ANALYSIS DATE : 06/17/06  
ANALYST : JIG

INSTRUMENT ID : GC/MS-02  
STD ID : PS061406-09

## VOLATILE ORGANIC COMPOUNDS BY EPA METHOD TO-14/TO-15

Continuing Calibration Verification of the 06/16/06 Calibration

| Compounds                    | Conc | Daily Conc | %REC |
|------------------------------|------|------------|------|
| BENZENE**                    | 10   | 10.61      | 106  |
| CARBON TETRACHLORIDE**       | 10   | 11.08      | 111  |
| Cyclohexane**                | 10   | 10.99      | 110  |
| 1,2-DICHLOROPROPANE**        | 10   | 10.52      | 105  |
| Bromodichloromethane**       | 10   | 10.91      | 109  |
| 1,4-Dioxane**                | 10   | 11.03      | 110  |
| TRICHLOROETHENE**            | 10   | 11.04      | 110  |
| 2,2,4-Trimethylpentane**     | 10   | 11.35      | 114  |
| Heptane**                    | 10   | 10.67      | 107  |
| cis- 1,3 DICHLOROPROPENE**   | 10   | 10.96      | 110  |
| MiBK**                       | 10   | 10.72      | 107  |
| trans 1,3 DICHLOROPROPENE**  | 10   | 11.17      | 112  |
| 1,1,2-TRICHLOROETHANE**      | 10   | 10.43      | 104  |
| TOLUENE**                    | 10   | 10.50      | 105  |
| 2-Hexanone**                 | 10   | 11.16      | 112  |
| Dibromochloromethane**       | 10   | 10.74      | 107  |
| 1,2 DIBROMOETHANE**          | 10   | 10.98      | 110  |
| TETRACHLOROETHYLENE**        | 10   | 10.27      | 103  |
| CHLOROBENZENE***             | 10   | 10.60      | 106  |
| ETHYLBENZENE***              | 10   | 10.64      | 106  |
| m-, & p- XYLENES***          | 20   | 21.57      | 108  |
| Bromoform***                 | 10   | 10.96      | 110  |
| STYRENE***                   | 10   | 10.86      | 109  |
| 1,1, 2,2- TETRACHLORETHANE** | 10   | 11.13      | 111  |
| o- XYLENE***                 | 10   | 11.00      | 110  |
| Ethyltoluene***              | 10   | 11.68      | 117  |
| 1,3,5- TRIMETHYLBENZENE***   | 10   | 11.51      | 115  |
| 1,2,4- TRIMETHYLBENZENE***   | 10   | 11.20      | 112  |
| Benzyl Chloride***           | 10   | 10.55      | 106  |
| 1,3- DICHLOROBENZENE***      | 10   | 11.00      | 110  |
| 1,4- DICHLOROBENZENE***      | 10   | 10.91      | 109  |
| 1,2-DICHLOROBENZENE***       | 10   | 11.22      | 112  |
| 1,2,4-TRICHLOROBENZENE***    | 10   | 9.32       | 93   |
| HEXACHLOROBUTADIENE***       | 10   | 11.16      | 112  |

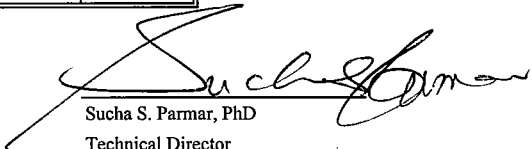
\* Internal std calculation IS1 : Bromochloromethane

\*\* Internal std calculation IS2 : 1,4-Difluorobenzene

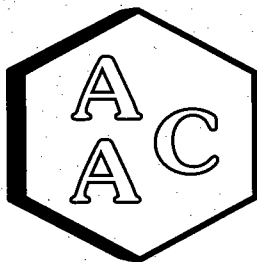
\*\*\* Internal std calculation IS3 : Chlorobenzene-d5

%REC should be 70-130%

!! Compound failed criteria and results should be considered estimated.

  
Sucha S. Parmar, PhD  
Technical Director





## Atmospheric Analysis & Consulting, Inc.

### Quality Control/Quality Assurance Report

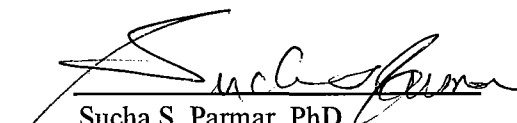
CLIENT ID : Laboratory Control Spike      DATE ANALYZED : 06/12/06  
AAC ID : LCS/LCSD      DATE REPORTED : 06/16/06  
MEDIA : Air      UNITS : ppbv

### TO-14/15 Laboratory Control Spike Recovery

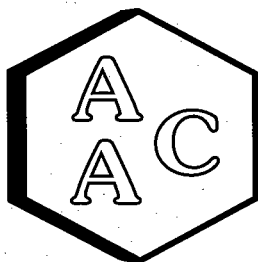
| Compound             | Sample Conc. | Spike Added | Spike Res | Dup Spike Res | Spike % Rec * | Spike Dup % Rec * | RPD** % |
|----------------------|--------------|-------------|-----------|---------------|---------------|-------------------|---------|
| 1,1-DICHLOROETHYLENE | 0.0          | 10.00       | 9.67      | 10.06         | 97            | 101               | 4.0     |
| METHYLENECHLORIDE    | 0.0          | 10.00       | 9.01      | 9.05          | 90            | 90                | 0.4     |
| BENZENE              | 0.0          | 10.00       | 8.88      | 9.23          | 89            | 92                | 3.9     |
| TRICHLOROETHENE      | 0.0          | 10.00       | 9.03      | 9.59          | 90            | 96                | 6.0     |
| TOLUENE              | 0.0          | 10.00       | 8.36      | 8.65          | 84            | 86                | 3.4     |
| TETRACHLOROETHYLENE  | 0.0          | 10.00       | 8.78      | 9.14          | 88            | 91                | 4.0     |
| CHLOROBENZENE        | 0.0          | 10.00       | 10.84     | 11.09         | 108           | 111               | 2.3     |
| ETHYLBENZENE         | 0.0          | 10.00       | 9.01      | 9.24          | 90            | 92                | 2.5     |
| m-, & p- XYLENES     | 0.0          | 20.00       | 18.21     | 18.74         | 91            | 94                | 2.9     |
| o- XYLENE            | 0.0          | 10.00       | 9.83      | 9.80          | 98            | 98                | 0.3     |

\* Must be 70-130%

\*\* Must be < 25%

  
Sucha S. Parmar, PhD  
Technical Director





# Atmospheric Analysis & Consulting, Inc.

ANALYSIS DATE : 06/12/06

INSTRUMENT ID : GC/MS-02

ANALYST : JJG

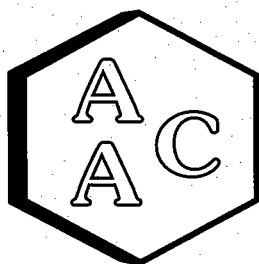
STD ID : PS052306-07

## VOLATILE ORGANIC COMPOUNDS BY EPA METHOD TO-14/TO-15

Continuing Calibration Verification of the 06/09/06 Calibration

| Compounds                      | Conc | Daily Conc | %REC |
|--------------------------------|------|------------|------|
| 4-BFB (surrogate standard)***  | 10   | 10.73      | 107  |
| Chlorodifluoromethane*         | 10   | 9.33       | 93   |
| Propylene*                     | 10   | 8.02       | 80   |
| DiCIDIFMethane*                | 10   | 9.75       | 98   |
| CHLOROMETHANE*                 | 10   | 9.50       | 95   |
| 1,2 DiCl-1,1,2,2-TetraFEthane* | 10   | 9.58       | 96   |
| VINYL CHLORIDE*                | 10   | 9.60       | 96   |
| Methanol*                      | 10   | 8.01       | 80   |
| 1,3-Butadiene*                 | 10   | 9.77       | 98   |
| BROMOMETHANE*                  | 10   | 9.51       | 95   |
| CHLOROETHANE*                  | 10   | 9.35       | 94   |
| Dichlorofluoromethane*         | 10   | 9.77       | 98   |
| Ethanol*                       | 10   | 8.91       | 89   |
| Vinyl Bromide*                 | 10   | 9.96       | 100  |
| Acetone*                       | 10   | 10.17      | 102  |
| TRICHLOROFLUOROMETHANE*        | 10   | 9.70       | 97   |
| Isopropanol*                   | 10   | 9.54       | 95   |
| Acrylonitrile*                 | 10   | 9.84       | 98   |
| 1,1 DICHLOROETHENE*            | 10   | 9.67       | 97   |
| METHYLENE CHLORIDE*            | 10   | 9.01       | 90   |
| Allyl CHLORIDE*                | 10   | 9.96       | 100  |
| Carbon disulfide*              | 10   | 9.58       | 96   |
| 1,1,2-TRICHLORO-1,2,2-TRIFLUO  | 10   | 9.42       | 94   |
| trans-1,2- DICHLOROETHYLENE*   | 10   | 9.25       | 93   |
| 1,1- DICHLOROETHANE*           | 10   | 9.44       | 94   |
| MTBE*                          | 10   | 9.81       | 98   |
| Vinyl Acetate*                 | 10   | 9.65       | 97   |
| MEK*                           | 10   | 9.85       | 99   |
| cis-1,2- DICHLOROETHYLENE*     | 10   | 11.11      | 111  |
| Hexane*                        | 10   | 11.06      | 111  |
| CHLOROFORM*                    | 10   | 11.37      | 114  |
| Ethyl Acetate*                 | 10   | 11.14      | 111  |
| Tetrahydrofuran*               | 10   | 11.17      | 112  |
| 1,2-DICHLOROETHANE*            | 10   | 11.02      | 110  |
| 1,1,1-TRICHLOROETHANE*         | 10   | 10.77      | 108  |





# Atmospheric Analysis & Consulting, Inc.

ANALYSIS DATE : 06/12/06  
ANALYST : JJG

INSTRUMENT ID : GC/MS-02  
STD ID : PS052306-07

## VOLATILE ORGANIC COMPOUNDS BY EPA METHOD TO-14/TO-15 Continuing Calibration Verification of the 06/09/06 Calibration

| Compounds                     | Conc | Daily Conc | %REC |
|-------------------------------|------|------------|------|
| BENZENE**                     | 10   | 8.88       | 89   |
| CARBON TETRACHLORIDE**        | 10   | 8.78       | 88   |
| Cyclohexane**                 | 10   | 8.92       | 89   |
| 1,2-DICHLOROPROPANE**         | 10   | 10.43      | 104  |
| Bromodichloromethane**        | 10   | 10.57      | 106  |
| 1,4-Dioxane**                 | 10   | 8.61       | 86   |
| TRICHLOROETHENE**             | 10   | 9.03       | 90   |
| 2,2,4-Trimethylpentane**      | 10   | 9.17       | 92   |
| Heptane**                     | 10   | 8.70       | 87   |
| cis- 1,3 DICHLOROPROPENE**    | 10   | 8.05       | 81   |
| MIBK**                        | 10   | 8.44       | 84   |
| trans 1,3 DICHLOROPROPENE**   | 10   | 8.14       | 81   |
| 1,1,2-TRICHLOROETHANE**       | 10   | 10.24      | 102  |
| TOLUENE**                     | 10   | 8.36       | 84   |
| 2-Hexanone**                  | 10   | 8.12       | 81   |
| Dibromochloromethane**        | 10   | 10.69      | 107  |
| 1,2 DIBROMOETHANE**           | 10   | 8.43       | 84   |
| TETRACHLOROETHYLENE**         | 10   | 8.78       | 88   |
| CHLOROBENZENE***              | 10   | 10.84      | 108  |
| ETHYLBENZENE***               | 10   | 9.01       | 90   |
| m-, & p- XYLENES***           | 20   | 18.21      | 91   |
| Bromoform***                  | 10   | 11.67      | 117  |
| STYRENE***                    | 10   | 8.59       | 86   |
| 1,1, 2,2- TETRACHLOROETHANE** | 10   | 11.23      | 112  |
| o- XYLENE***                  | 10   | 9.83       | 98   |
| Ethyltoluene***               | 10   | 9.81       | 98   |
| 1,3,5- TRIMETHYLBENZENE***    | 10   | 10.34      | 103  |
| 1,2,4- TRIMETHYLBENZENE***    | 10   | 10.27      | 103  |
| Benzyl Chloride***            | 10   | 8.41       | 84   |
| 1,3- DICHLOROBENZENE***       | 10   | 10.02      | 100  |
| 1,4- DICHLOROBENZENE***       | 10   | 9.46       | 95   |
| 1,2-DICHLOROBENZENE***        | 10   | 9.89       | 99   |
| 1,2,4-TRICHLOROBENZENE***     | 10   | 8.74       | 87   |
| HEXACHLOROBUTADIENE***        | 10   | 11.44      | 114  |

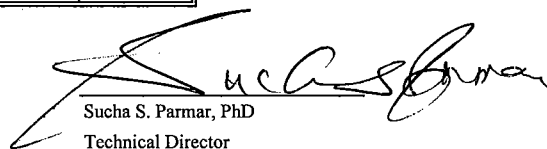
\* Internal std calculation IS1 : Bromochloromethane

\*\* Internal std calculation IS2 : 1,4-Difluorobenzene

\*\*\* Internal std calculation IS3 : Chlorobenzene-d5

%REC should be 70-130%

!! Compound failed criteria and results should be considered estimated.

  
Sucha S. Parmar, PhD  
Technical Director



060263

AAC

**CLIENT**

Redwood Landfill

DATE: 6-8-06

PAGE: 1 of 1

[illegible]

***Appendix E***  
***Source Test Plan and***  
***California Air Resources Board Independent***  
***Contractor Certification***



Shaw Environmental, Inc.

2360 Bering Drive  
San Jose, California 95131  
408-382-5800  
FAX: 408-433-1912

**VIA FAX 415-749-4922**

To: Tim Underwood, BAAQMD  
Subject: Source Test Protocol & Test Notification

Date: May 24, 2006

**SOURCE TEST PROTOCOL  
FOR  
REDWOOD LANDFILL  
FLARES, A-50 & A-51**

Shaw Environmental, Inc. (Shaw) will perform compliance testing on landfill gas flares A-50 and A-51 located at the Redwood Landfill in Novato, California. The purpose of the test is to demonstrate the performance of the landfill gas flare as specified by the Bay Area Air Quality Management District (BAAQMD) Major Facility Review (MFR) Title V permits and Regulation 8, Rule 34. This protocol is to inform the BAAQMD of the planned test date and testing procedures. Shaw is approved by the California Air Resources Board (ARB) as an independent contractor to conduct compliance emission testing.

**Source Test Information, Redwood Landfill**

|                          |                                                                                                                                                  |                                                                                                                                                                                         |
|--------------------------|--------------------------------------------------------------------------------------------------------------------------------------------------|-----------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------|
| Test Location:           | Redwood Landfill, Inc.<br>8950 Redwood Highway<br>Novato, CA 94948<br>BAAQMD MFR Number A1179                                                    | Contact: Ms. Beth Shiverdecker<br>Telephone: 925-719-0016                                                                                                                               |
| Source to be Tested:     | Flares A-50 and A-51                                                                                                                             |                                                                                                                                                                                         |
| Test Objective:          | Determine compliance with Title V Permit and Regulation 8, Rule 34<br>Flare: Permit Condition Number 19867                                       | Emission Limits:<br>NO <sub>x</sub> : 15 ppm @3% O <sub>2</sub><br>CO: 123 ppmv @3% O <sub>2</sub><br>NMOC: > 98% destruction efficiency<br>or < 30 ppm as methane at 3% O <sub>2</sub> |
| Test to be Performed By: | Shaw Environmental, Inc.<br>2360 Bering Drive<br>San Jose, CA 95131                                                                              | Contact: Bill Johnston<br>Tel: (408) 382-5822                                                                                                                                           |
| Test date                | June 7&8, 2006                                                                                                                                   |                                                                                                                                                                                         |
| Test Parameters:         | Inlet<br>O <sub>2</sub> , CO <sub>2</sub> , N <sub>2</sub> , HHV<br>THC, CH <sub>4</sub> , NMOC, Sulfur Species,<br>VOCs, landfill gas flow rate | Outlet<br>NMOC, NO <sub>x</sub> , CO,<br>Volumetric flow rate                                                                                                                           |

## SCOPE OF WORK

A source test will be performed on flares A-50 and A-51 located at the Redwood Landfill. The flare source test shall determine the following parameters as specified in the facility's Title V permit condition number 19867 and BAAQMD Regulation 8, Rule 34. The following items in the condition will be met.

- Determine landfill gas flow rate (dry basis);
- The landfill gas shall be analyzed for higher heating value (HHV), carbon dioxide (CO<sub>2</sub>), nitrogen (N<sub>2</sub>), oxygen (O<sub>2</sub>), total hydrocarbons (THC), methane (CH<sub>4</sub>), non-methane organic compounds (NMOC), speciated organic and sulfur compounds. All concentrations shall be reported on a dry basis;
- Stack gas flow rate from the flare;
- Concentrations (dry basis) of NMOC, CO, NO<sub>x</sub>, and O<sub>2</sub> in the flare stack gas;
- NO<sub>x</sub> and CO emission rates will be reported in units of lbs/MMBTU;
- NMOC destruction efficiency achieved by the flare, if the outlet NMOC is greater than 30 parts per million by volume (ppmv), expressed as methane at 3 percent oxygen on a dry basis.;
- The average combustion temperature in the flare during the test period.
- A characterization of the landfill gas for all the compounds listed below, as identified in Part 31-a of the condition.

The source test will demonstrate the following limits.

- Emissions of NO<sub>x</sub> shall not exceed 0.06 pounds per million BTU (lb/MMBtu) or exceed 15 ppmv corrected to 15 percent O<sub>2</sub>.
- Emissions of CO shall not exceed 0.30 lb/MMBtu or exceed 123 ppmv corrected to 15 percent O<sub>2</sub>.
- The flare destruction efficiency of NMOC shall not be less than 98 percent by weight, unless the outlet NMOC concentration is less than 30 ppmv, expressed as methane at 3 percent oxygen on a dry basis.
- Concentrations of selected organic and sulfur compounds, as identified in Part 18, a and b of the condition.

## TEST PROCEDURES

The source test matrix identified below presents the test parameters and source test methods (each compliant with the requirements of the BAAQMD) that will be utilized to determine the emissions from the flare.

### Landfill Gas Fuel Analysis

Integrated Tedlar bag samples will be collected at the inlet to the flare. Samples will be collected concurrently with the outlet sampling. Samples will be analyzed within 72 hours (48 hrs for sulfur). One of the samples will be analyzed for speciated organics and sulfur compounds.

### Continuous Emission Monitoring

Single point sampling will be performed at the center of the stack. All analyzers will be checked for calibration before and after each run. Emission monitoring data will be recorded on strip charts and a data logger. Methane concentrations will be determined by passing sample gas through an activated carbon scrubber prior to the hydrocarbon analyzer. Hydrocarbon emissions will be measured and reported as methane.

## Process Parameters

Flare process parameters including the fuel gas flow rate and the flare operating temperature will be measured and recorded during each run using the flare gas measurement system.

### Test Matrix Redwood Landfill LFG and Flares A-50 & A-51

| Sample Parameter                        | Test Method               | Runs | Test Duration |
|-----------------------------------------|---------------------------|------|---------------|
| <b>LFG Parameters- Flare Inlet</b>      |                           |      |               |
| Gas flow rate                           | Flow meter                | 3    | 30 Minutes    |
| Fuel factor                             | EPA Method 19             | 3    | 30 Minutes    |
| Fixed gases <sup>1</sup>                | ASTM D1945, EPA Method 18 | 1    |               |
| Volatile Organic Compounds <sup>2</sup> | EPA Method TO 15          | 1    | 30 Minutes    |
| Sulfur compounds <sup>2</sup>           | ASTM D5504                | 1    | 30 Minutes    |
| NMOC                                    | EPA Method 25 C           | 1    | 30 Minutes    |
| <b>Source Parameters-Flare Exhaust</b>  |                           |      |               |
| Volumetric flow rate                    | EPA Method 19             | 3    | 30 Minutes    |
| O <sub>2</sub>                          | BAAQMD ST-14              | 3    | 30 Minutes    |
| NMOC                                    | BAAQMD ST-7               | 3    | 30 Minutes    |
| CO                                      | BAAQMD ST-6               | 3    | 30 Minutes    |
| NOx                                     | BAAQMD ST-13A             | 3    | 30 Minutes    |

<sup>1</sup> Fixed gases include HHV, CO<sub>2</sub>, N<sub>2</sub>, O<sub>2</sub>, THC and CH<sub>4</sub>.

<sup>2</sup> Only one sample will be analyzed for the test program.

## REPORT

The final report is due to the BAAQMD 45 days after testing has occurred. The report will describe the tests that were conducted, the operating conditions of the source during the test, and the emission results reported in units of the appropriate standard. All raw data and sample calculations used to obtain the reported results will be included so that the accuracy of the reported results can be verified from the report. The flare process information will be submitted in the report.

## TEST SCHEDULE

The source test is scheduled for June 7&8, 2006. One flare will be tested per day.

State of California  
Air Resources Board  
Approved Independent Contractor  
**Shaw Environmental, Inc.**

This is to certify that the company listed above has been approved  
by the Air Resources Board to conduct compliance testing  
pursuant to section 91207, title 17, California Code of Regulations,  
until June 30, 2006 for those test methods listed below:

ARB Source Test Methods:  
1, 2, 3, 4, 5, 8, 100 (CO, CO<sub>2</sub>, NO<sub>x</sub>, O<sub>2</sub>, SO<sub>2</sub>, THC)  
Visible Emissions Evaluation

  
William V. Loscutoff, Chief  
Monitoring and Laboratory Division

***Appendix F***  
***Major Facility Review Permit***

# **Bay Area Air Quality Management District**

939 Ellis Street  
San Francisco, CA 94109  
(415) 771-6000

---

**Final**

## **MAJOR FACILITY REVIEW PERMIT**

**Issued To:**  
**Redwood Landfill, Inc.**  
**Facility #A1179**

**Facility Address:**  
8950 Redwood Highway  
Novato, CA 94948

**Mailing Address:**  
P. O. Box 793  
Novato, CA 94948

**Responsible Official**  
Ramin Khany  
Landfill Manager  
415-892-2851

**Facility Contact**  
Beth Shiverdecker  
Environmental Compliance Specialist  
415-892-2851

---

|                          |                                   |                                             |
|--------------------------|-----------------------------------|---------------------------------------------|
| <b>Type of Facility:</b> | Landfill for Solid Waste Disposal | <b>BAAQMD Engineering Division Contact:</b> |
| <b>Primary SIC:</b>      | 4953                              | Carol S. Allen                              |
| <b>Product:</b>          | Refuse and Sludge Disposal        |                                             |

**ISSUED BY THE BAY AREA AIR QUALITY MANAGEMENT DISTRICT**

Signed by Jack P. Broadbent  
Jack P. Broadbent, Executive Officer/Air Pollution Control Officer

April 18, 2006  
Date

## VI. Permit Conditions

### Condition # 19867

**FOR: S-5 REDWOOD LANDFILL WITH GAS COLLECTION SYSTEM; A-18 WATER SPRAYS; A-50 LANDFILL GAS FLARE; AND A-51 LANDFILL GAS FLARE**

- |                                        | Required Components |
|----------------------------------------|---------------------|
| Total Number of Vertical Wells:        | 86                  |
| Total Number of Horizontal Collectors: | 8                   |
- b. The Permit Holder has been issued an Authority to Construct for the landfill gas collection system components listed below. Specific well and collector locations, depths, and lengths of associated piping are as described in detail in Permit Application # 13027.

|                                    | Minimum | Maximum |
|------------------------------------|---------|---------|
| Install New Vertical Wells:        | 0       | 30      |
| Decommission Vertical Wells:       | 0       | 18      |
| Install New Horizontal Collectors  | 0       | 10      |
| Decommission Horizontal Collectors | 0       | 4       |
| Replace Vertical Wells *           | 0       | 15      |

\* one-for-one well replacement at new optimal locations

Wells installed or shutdown pursuant to subpart b shall be added to or removed from subpart a in accordance with the procedures identified in Regulations 2-6-414 or 2-6-415. The Permit Holder shall maintain records of the decommissioning date for each well that is shut down and the initial operation date for each new well.

(Basis: Regulations 2-1-301, 8-34-301.1, 8-34-304, 8-34-305, and 2-6-413)

18. If a gas characterization test indicates that this site's landfill gas contains organic compounds in excess of any of the concentrations listed in Parts 18a or 18b below, then the Permit Holder shall submit an application for a Change of Permit Conditions, within no later than 30 days from receipt of the test results.

- a. Total Non-Methane Organic Compounds: 750 ppmv  
 (calculated as hexane equivalent)  
 (Basis: Cumulative Increase)

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\*b. For toxic air contaminants (TACs):

| <u>Compound</u>           | <u>Concentration</u> |
|---------------------------|----------------------|
| Acrylonitrile             | 280 ppbv             |
| Benzene                   | 340 ppbv             |
| Carbon Tetrachloride      | 70 ppbv              |
| Chloroform                | 70 ppbv              |
| 1,4 Dichlorobenzene       | 400 ppbv             |
| 1,1 Dichloroethane        | 150 ppbv             |
| Ethylene Dibromide        | 70 ppbv              |
| Ethylene Dichloride       | 70 ppbv              |
| Methylene Chloride        | 320 ppbv             |
| Perchloroethylene         | 450 ppbv             |
| 1,1,2,2 Tetrachloroethane | 70 ppbv              |
| Trichloroethylene         | 250 ppbv             |
| Vinyl Chloride            | 880 ppbv             |

(Basis: Toxic Risk Management Policy)

c. The concentration of total reduced sulfur compounds (TRS) in collected landfill gas shall not exceed a peak of 1300 ppmv (calculated as H<sub>2</sub>S) and shall not exceed an annual average of 425 ppmv (calculated as H<sub>2</sub>S). The peak and annual average TRS concentrations shall be measured and calculated in accordance with Parts 31a and 31b.

(Basis: Cumulative Increase, RACT, and Regulation 9-1-302)

19. The A-50 and A-51 Landfill Gas Flares shall be fired on landfill gas and may also be used to abate leachate vapors from the S-50 Leachate Vaporator. (Basis: RACT and Regulation 2-2-112)
20. The total combined throughput of landfill gas (with an HHV of 500 BTU/scf) to the A-50 Landfill Gas Flare and the A-51 Landfill Gas Flare shall not exceed 1,490,000,000 scf during any consecutive 12-month period and shall not exceed 5,760,000 scf during any one day. In order to demonstrate compliance with this condition, the A-50 and A-51 Flares shall be equipped with a one or more properly operating continuous gas flow meters. (Basis: Cumulative Increase, 40 CFR 60.756(b)(2)(i))
21. [deleted]

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22. The temperature in the combustion zone of each flare shall be maintained at the minimum temperature listed below, averaged over any 3-hour period. In order to demonstrate compliance with this condition, A-50 and A-51 shall each be equipped with a continuous temperature monitor and recorder. If a source test demonstrates compliance with all applicable requirements at a different temperature, the APCO may revise these temperature limits, in accordance with the procedures identified in Regulation 2-6-414 or 2-6-415, based on the following criteria. The minimum combustion zone temperature for the flare shall be equal to the average combustion zone temperature determined during the most recent complying source test minus 50 degrees F, provided that the minimum combustion zone temperature is not less than 1400 degrees F.
- a. The minimum combustion zone temperature for A-50 is 1475 degrees F, averaged over any 3-hour period.
  - b. The minimum combustion zone temperature for A-51 is 1422 degrees F, averaged over any 3-hour period.
- (Basis: Toxic Risk Management Policy, Regulations 8-34-301.3 and 8-34-501.3, and 40 CFR 60.756(b)(1))
23. The A-50 and A-51 Landfill Gas Flares shall comply with the NMOC emission limit in Regulation 8-34-301.3.  
(Basis: Cumulative Increase, 8-34-301.3, and 40 CFR 60.752(b)(2)(iii)(B))
- \*24. The A-50 and A-51 Landfill Gas Flares shall each achieve a minimum destruction efficiency of 83% by weight for any EPA Hazardous Air Pollutants or any District toxic compounds that are determined to be present in the landfill gas or leachate vapors. (Basis: Toxic Risk Management Policy)
25. Nitrogen oxides (NO<sub>x</sub>) emissions from each enclosed flare (A-50 and A-51) shall not exceed 0.06 pounds of NO<sub>x</sub>, calculated as NO<sub>2</sub>, per million BTU. Compliance with this emission limit may be demonstrated by not exceeding the following flue gas concentration limit: 15 ppmv of NO<sub>x</sub>, corrected to 15% oxygen, dry basis.  
(Basis: RACT and Offsets)
26. Carbon monoxide (CO) emissions from each enclosed flare (A-50 and A-51) shall not exceed 0.30 pounds of CO per million BTU. Compliance with this emission limit may be demonstrated by not exceeding the following flue gas concentration limit: 123 ppmv of CO, corrected to 15% oxygen, dry basis.  
(Basis: RACT and Cumulative Increase)

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27. [deleted]
28. [deleted]
29. The Permit Holder shall maintain records of all planned and unanticipated shut downs of the A-50 and A-51 Flares and of any temperature excursions. The records shall include the date, time, duration, and reason for any shut down or excursion. Any unanticipated shut downs or temperature excursions shall be reported to the Enforcement Division immediately. All inspection and maintenance records, records of shut downs and excursions, gas flow records, temperature records, analytical results, source test results, and any other records required to demonstrate compliance with the above permit conditions, Regulation 8 Rule 34, or 40 CFR Part 60 Subpart WWW shall be retained on site for a minimum of five years and shall be made available to District staff upon request. (Basis: 2-6-501, 8-34-501, 40 CFR 60.758)
30. In order to demonstrate compliance with Parts 23, 25, and 26 above, Regulation 8, Rule 34, Sections 301.3 and 412, and 40 CFR 60.8 and 60.752(b)(2)(iii)(B), the Permit Holder shall ensure that a District approved source test is conducted annually on the A-50 Landfill Gas Flare and the A-51 Landfill Gas Flare. Each annual source test shall determine the following:
  - a. landfill gas flow rate to the flare (dry basis);
  - b. concentrations (dry basis) of carbon dioxide (CO<sub>2</sub>), nitrogen (N<sub>2</sub>), oxygen (O<sub>2</sub>), total hydrocarbons (THC), methane (CH<sub>4</sub>), and total non-methane organic compounds (NMOC) in the landfill gas;
  - c. stack gas flow rate from the flare (dry basis);
  - d. concentrations (dry basis) of NO<sub>x</sub>, CO, NMOC, and O<sub>2</sub> in the flare stack gas;
  - e. NMOC destruction efficiency achieved by the flare;
  - f. NO<sub>x</sub> and CO emission rates from the flare in units of pounds per MM BTU,
  - g. average combustion zone temperature in the flare during the test period.

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The first source test for A-51 shall be conducted no later than 60 days after the initial start-up date for A-51. Each annual source test shall be conducted no earlier than 9 months and no later than 12 months after the previous annual source test. The Source Test Section of the District shall be contacted to obtain approval of the source test procedures at least 14 days in advance of each source test. The Source Test Section shall be notified of the scheduled test date at least 7 days in advance of each source test. The source test report shall be submitted to the Compliance and Enforcement Division and the Source Test Section within 60 days of the test date.

(Basis: Cumulative Increase, Toxic Risk Management Policy, RACT, Offsets, Regulations 8-34-301.3, 8-34-412, 40 CFR 60.8 and 40 CFR 60.752(b)(2)(iii)(B))

31. Landfill Gas Testing:

- a. The Permit Holder shall conduct a characterization of the landfill gas on a quarterly basis with one test concurrent with one of the annual source tests required by Part 30 above. The landfill gas sample shall be drawn from the main landfill gas header. Each quarterly landfill gas sample shall be analyzed for the sulfur compounds listed below. Once per year (concurrent with a Part 30 annual source test) the landfill gas shall be analyzed for all the organic and sulfur compounds listed below. All concentrations shall be reported on a dry basis. The laboratory analysis report for the annual organic and sulfur compound gas characterization test shall be included with the Part 30 source test report and shall be submitted to the Compliance and Enforcement Division and the Source Test Section within 60 days of the test date. (Basis: Toxic Risk Management Policy and Regulations 8-34-412 and 9-1-302)

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| <u>Organic Compounds</u> | <u>Organic Compounds</u>  | <u>Sulfur Compounds</u> |
|--------------------------|---------------------------|-------------------------|
| acrylonitrile            | ethylene dibromide        | carbon disulfide        |
| benzene                  | fluorotrichloromethane    | carbonyl sulfide        |
| carbon tetrachloride     | hexane                    | dimethyl sulfide        |
| chlorobenzene            | isopropyl alcohol         | ethyl mercaptan         |
| chlorodifluoromethane    | methyl ethyl ketone       | hydrogen sulfide        |
| chloroethane             | methylene chloride        | methyl mercaptan        |
| chloroform               | perchloroethylene         |                         |
| 1,1 dichloroethane       | toluene                   |                         |
| 1,1 dichlorethene        | 1,1,1 trichloroethane     |                         |
| 1,2 dichlorethene        | 1,1,2,2 tetrachloroethane |                         |
| 1,4 dichlorbenzene       | trichloroethylene         |                         |
| dichlorodifluoromethane  | vinyl chloride            |                         |
| dichlorofluoromethane    | xylene                    |                         |
| ethylbenzene             |                           |                         |

- b. Once per week, beginning no later than March 31, 2005, the Permit Holder shall analyze the landfill gas for hydrogen sulfide (H<sub>2</sub>S) concentration using a Draeger tube to further demonstrate compliance with Part 18c and Regulation 9-1-302. The landfill gas sample shall be drawn from the main landfill gas header. The Permit Holder shall follow the manufacturer's procedures for using the Draeger tube and interpreting the results. The total reduced sulfur (TRS) content of the landfill gas shall be calculated using the average ratio of TRS/H<sub>2</sub>S for this site according to the following equation:  $TRS = 1.015 * H_2S \text{ measured by Draeger tube}$ . The Permit Holder shall maintain records of all Draeger tube test dates and test results and shall summarize the average H<sub>2</sub>S concentrations and the calculated TRS content of the landfill gas on a quarterly basis. Each Draeger tube test result (after conversion to TRS content) and the quarterly laboratory analysis in Part 31a shall be compared to the Peak TRS Limit in Part 18c. On a rolling quarterly basis, the Permit Holder shall determine the annual average TRS content for comparison to the Annual Average TRS Limit in Part 18c. (Basis: Cumulative Increase, RACT, and Regulation 9-1-302).

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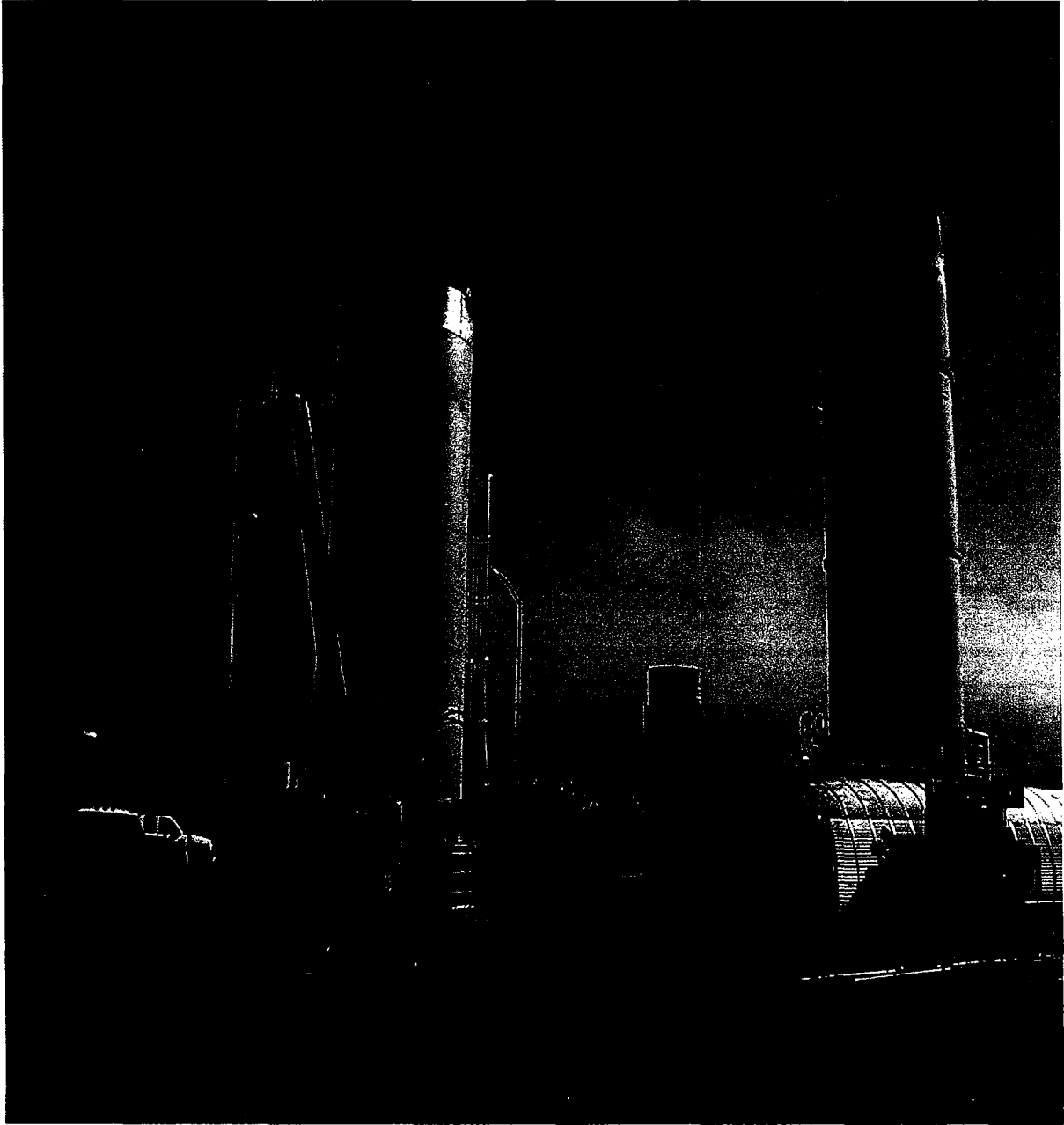
### **Condition # 19867**

**FOR: S-5 REDWOOD LANDFILL WITH GAS COLLECTION SYSTEM; A-18 WATER  
SPRAYS; A-50 LANDFILL GAS FLARE; AND A-51 LANDFILL GAS FLARE**

32. The annual report required by BAAQMD Regulation 8-34-411 shall be submitted in two semi-annual increments. The reporting period for the first increment of the Regulation 8-34-411 annual report that is submitted subsequent to the issuance of the MFR Permit for this site shall be from December 1, 2003 through April 30, 2004. This first increment report shall be submitted by May 31, 2004. The reporting periods and report submittal due dates for all subsequent increments of the Regulation 8-34-411 report shall be synchronized with the reporting periods and report submittal due dates for the semi-annual MFR Permit monitoring reports that are required by Section I.F of the MFR Permit for this site. A single report may be submitted to satisfy the requirements of Section I.F, Regulation 8-34-411, and 40 CFR Part 63.1980(a), provided that all items required by each applicable reporting requirement are included in the single report.  
(Basis: Regulation 8-34-411 and 40 CFR Part 63.1980(a))

***Appendix G***  
***Site Picture***

## Redwood Landfill Flares



A-51

A-50