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# **Emissions from Shell Core Making** and Storage

1413-115 HO

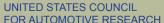
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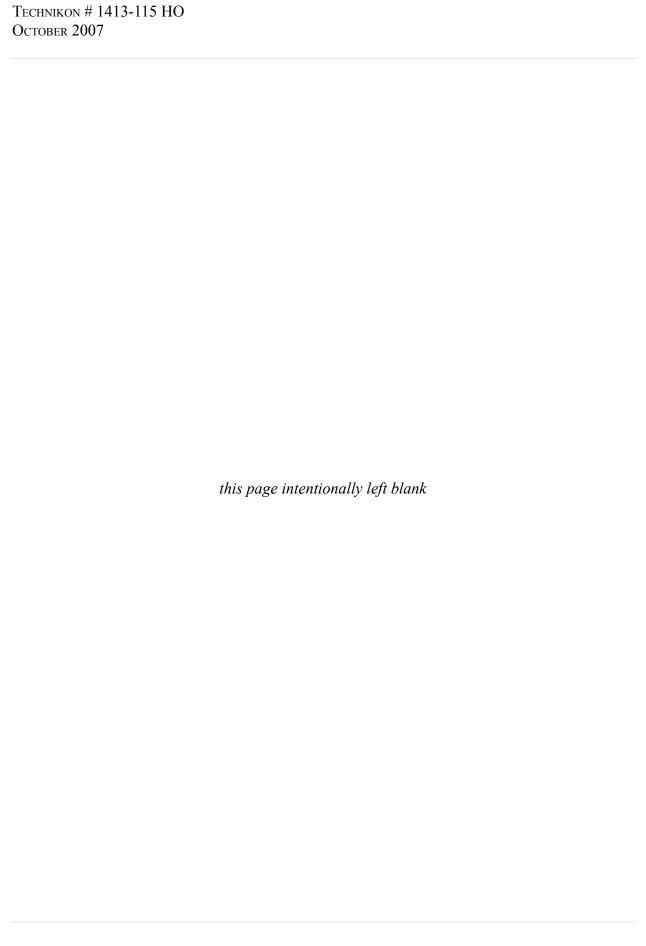








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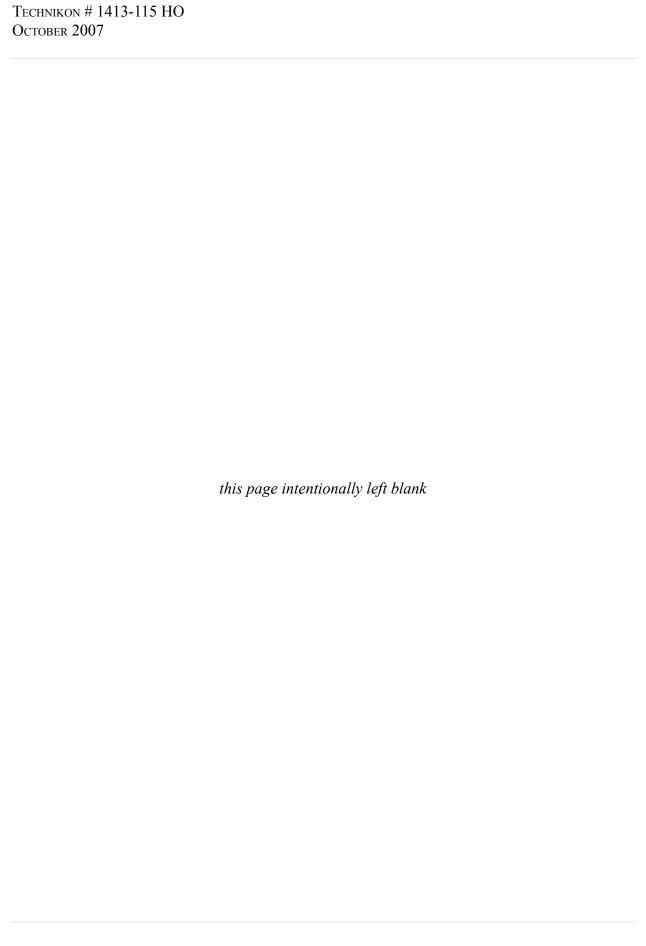
# **Emissions from Shell Core Making and Storage**

# 1413-115 HO

This report has been reviewed for completeness and accuracy and approved for release by the following:

Director of Measurement Technologies	//Original Signed//	
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Vice President	//Original Signed//	
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The data contained in this report were developed to assess the relative emissions profile of the product or process being evaluated. You may not obtain the same results in your facility. Data were not collected to assess cost or producibility.



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# **Executive Summary**

A testing protocol was undertaken for Test HO to quantitatively evaluate airborne emissions from shell core making and storage. A phenol-formaldehyde resin coated sand (Technisand<sup>TM</sup> 630BN Gold, Fairmount Minerals), which does not contain hexamethylenetetramine, was the core making material used for the test. The stated concentration of the binder was 3% based on sand (BOS) weight. Results were compared to Test HN, a baseline shell core making and storage test which used a hexamine containing resin (Technisand<sup>TM</sup> XC30 Shell Sand, Fairmount Minerals).

A Beardsley and Piper Cormatic Shell core machine, model SF6CA, with 32 natural gas burners, circa 1963, was used for the test. An insulated vertical draft hood draped on three sides with fire resistant and inert fabric and enclosed on the fourth with a wall of aluminum sheeting functioned as the shell core machine enclosure for capturing emissions. The shell machine was configured with the step-core pattern corebox. Emissions were measured from six replicate runs consisting of the making and concurrent storage of 15 cores, immediately followed by a separate storage period within the enclosure after the blow and cure cycle of the final core, to complete an elapsed run time of 60 minutes. Emission samples were continuously collected for the entire core making/storage and separate storage periods.

The emission results were calculated in both pounds of pollutant per pound of binder (lb/ lb) and pounds of pollutant per ton of cores (lb/ton). All emissions have been corrected by subtracting background contributions. Subtracting those compounds present from quality control and method background blanks (i.e. ambient background) generates emissions for the complete shell core making process, including machine, natural gas burner, and material emissions. Material (binder) emissions were obtained by subtracting the background contribution produced from a system blank obtained by operating the shell core machine without actual core production. Subtracting the system blank allows comparisons to be made for emissions between material systems only, without contributions from the shell core machine.

The system background contained moderate concentrations of numerous hydrocarbons in-

cluding benzene, toluene, phenol, and xylenes. The measured criteria pollutants and greenhouse gases concentration were also high in the blank. Comparing system blank concentrations to sample concentrations shows that a majority of the differentiated hydrocarbons, the total hydrocarbons as measured by TGOC as Propane, and the methane originated from the shell core machine, including the natural gas burners. It is also apparent that most of the total hydrocarbon emissions as measured by TGOC as Propane is comprised of methane. The emissions for total hydrocarbons and methane parallel each other exactly over the entire core making/storage and storage run.

Total hydrocarbon results for both lb/ton metal and lb/lb binder for Test HO show a statistically significant increase of 75% when compared to Test HN, as shown in Table 1a and Table 1b. Relative decreases of approximately 70% were shown for the other Emission Indicators including the Sum of Target Analytes and the Sum of Target HAPs. The difference between the results of the two tests for non-methane hydrocarbons was found to be insignificant.

Table 1a Average Emission Indicators Summary Table, Core Make/Store and Storage, Test HO to Test HN, Ib/tn Core

			Percent Change
Analyte Name	Reference Test HN	Test HO	from Reference
THC as Propane	5.34E-01	9.34E-01	75
Non-Methane Hydrocarbons	3.28E-01	2.96E-01	-10
Sum of Target Analytes	5.79E-01	1.46E-01	-75
Sum of Target HAPs	1.25E-01	4.06E-02	-67
Sum of Target POMs	ND	4.69E-05	NA

ND=Not Detected NA=Not Applicable

Table 1b Average Emission Indicators Summary Table, Core Make/Store and Storage, Test HO to Test HN, lb/lb Binder

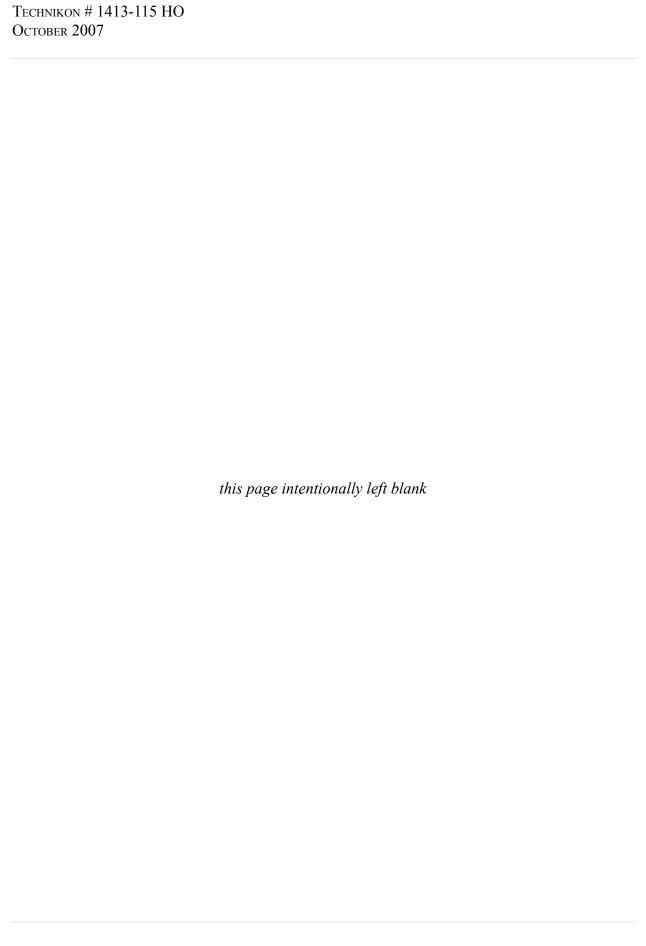
			Percent Change
Analyte Name	Reference Test HN	Test HO	from Reference
THC as Propane	9.16E-03	1.60E-02	75
Non-Methane Hydrocarbons	5.63E-03	5.08E-03	-10
Sum of Target Analytes	1.00E-02	2.53E-03	-75
Sum of Target HAPs	2.17E-03	7.07E-04	-67
Sum of Target POMs	ND	8.15E-07	NA

ND=Not Detected NA=Not Applicable Two target analytes comprised over 97% of emissions for the system background adjusted data. Ammonia had the highest contribution at 78% followed by phenol at 19%. Formaldehyde accounted for approximately 2% of emissions. The few compounds which supplied the remainder of measurable emissions were all less than 0.1%. Targeted HAPs contributed approximately 22% to total measured emissions.

Five target analyte emissions were lower and two were higher when comparing the two binder systems from Test HO and Test HN. The largest emissions decrease of 77% was found for ammonia, an EPA EPCRA Section 313 chemical, although its contribution (78%) was exactly the same as in Test HN. The largest increase in emissions of 81% was for crotonaldehyde.

A GC/MS analysis of the stack emissions revealed several additional compounds for both Test HO and Test HN. Some of these compounds were of higher concentration than the predicted compounds. Tentative identification included several silicone compounds (from the release agent) and numerous nitrogen containing compounds, including nitro phenols.

Emission results from the testing performed and described herein are not suitable for use as emission factors or for purposes other than evaluating the relative emission reductions associated with the use of alternative materials, equipment, or processes. The emissions measurements are unique to the specific cores produced, materials used, and testing methodology associated with these tests. These measurements should not be used as the basis for estimating emissions from actual commercial foundry applications.



### 1.0 Introduction

# 1.1. Background

Technikon LLC is a privately held contract research organization located in McClellan, California, a suburb of Sacramento. Technikon offers emissions research services to industrial and government clients specializing in the metal casting and point source emissions areas. Technikon operates the Casting Emission Reduction Program (CERP). CERP is a cooperative initiative between the Department of Defense (US Army) and the United States Council for Automotive Research (USCAR). The parties to the CERP Cooperative Research and Development Agreement (CRADA) include The Environmental Leadership Council of USCAR, a Michigan partnership of DaimlerChrysler Corporation, Ford Motor Company, and General Motors Corporation; the U.S. Army Research, Development, and Engineering Command (RDECOM-ARDEC); the American Foundry Society (AFS); and the Casting Industry Suppliers Association (CISA). The US Environmental Protection Agency (US EPA) and the California Air Resources Board (CARB) also have been participants in the CERP program and rely on CERP published reports for regulatory compliance data. All published reports are available on the CERP web site at <a href="https://www.cerp-us.org">www.cerp-us.org</a>.

### 1.2. CERP/Technikon Objectives

The primary objective of CERP is to evaluate materials, equipment, and processes used in the production of metal castings. Technikon's facility was designed to evaluate alternative materials and production processes while achieving significant air emission reductions. The facility's principal testing arena is designed to measure airborne emissions from individually poured molds. This testing facility enables the repeatable collection and evaluation of airborne emissions and associated process data.

# 1.3. Report Organization

This report has been written to document the methodology and results of a specific test plan

that was used to evaluate airborne emissions from shell core making and storage. Shell cores were made from resin coated sand containing 3% binder BOS.

Section 2.0 of this report includes a summary of the methodologies used for data collection and analysis, procedures for emission calculations, QA/QC procedures, and data management and reduction methods. Specific data collected during this test are summarized in Section 3.0 and detailed data appear in the appendices of this report. Section 4.0 contains a discussion of the results.

The raw data for this test series are archived at the Technikon facility.

# 1.4. Specific Test Plan and Objectives

Test HO was designed and conducted to evaluate airborne emissions from the making and storage of shell cores. Shell step cores were made with Technisand<sup>TM</sup> 630BN Gold Shell Sand from Fairmount Minerals containing 3% resin BOS.

Table 1-1 provides a summary of the test plan. The details of the approved test plan are included in Appendix A.

Table 1-1 Test Plan Summary

Type of Process Tested	Shell Core Making	Shell Core Making
Test Plan Number	1413-122 HN	1413-115 HO
Core	3 % (BOS) Fairmount Minerals XC30 Shell sand	3 % (BOS) Fairmount Minerals 630BN Gold Shell sand
Number of Runs	6 runs with 15 cores made	6 runs with 15 cores made
Test Dates	April 23, 2007 through April 25, 2007	April 25, 2007 through April 26, 2007
Emissions Measured	81 target analytes and TGOC as Propane, CO, CO <sub>2</sub> , NOx, SO <sub>2</sub> , CH <sub>4</sub> Propane, CO, CO2, NOx, SO <sub>2</sub>	
Process Parameters Measured	Core weight, binder weight, core LOI, core box temperature	Core weight, binder weight, core LOI, core box temperature

#### 2.0 Test Methods and Procedures

# 2.1. Testing Program

The testing program encompasses the foundry process and emissions testing, both of which are rigorously controlled. Relevant parameters are monitored and recorded prior to and during the emission tests. Process measurements for the shell core tests included the weights of the cores, loss on ignition (LOI) values for the core prior to the test, and the temperature of the corebox. Measured source parameters included stack temperature, pressure, volumetric flow rate, and moisture content. All parameters were maintained within prescribed ranges to ensure the reproducibility of the test runs.

# 2.1.1. Test Plan Review and Approval

The proposed test plan was reviewed and approved by the Technikon staff and by CERP Working Group Chairs as appropriate. The Test Plans for Test HO and the baseline shell core Test HN are included in Appendix A.

# 2.1.2. Shell Core Making and Storage

Figure 2-1 is a diagram of the shell core making and storage test setup.

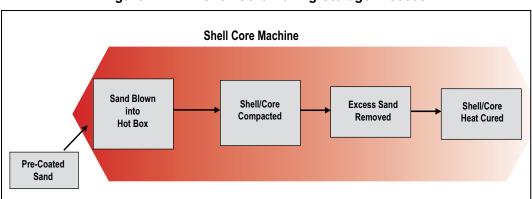


Figure 2-1 Shell Core Making/Storage Process

Table 2-1 Beardsley and Piper SF Cormatic Specifications

Sand magazine capacity	60 lbs
Core box: Maximum height	20 in
Maximum length	15.5 in
Maximum width	21 in
BTU per hour natural gas	90000
Voltage	125 VAC
Amps	8

Figure 2-2 Shell Core Machine Enclosure



Figure 2-3 Cross Section of Shell Step Core



Shell cores were made on a standard vertically parted semi-automated machine. A Beardsley and Piper Cormatic Shell core machine, model SF6CA, with 32 natural gas burners, circa 1963 was used for the test. Machine specifications are summarized in Table 2-1. The shell machine was configured with the step core corebox. This pattern was specifically built to evaluate core emissions. An insulated vertical draft hood draped on three sides with fire resistant and chemically inert fabric and enclosed on the fourth side with a wall of aluminum sheeting functioned as the shell core machine enclosure for capturing emissions. The machine within the enclosure is shown in Figure 2-2.

Prior to the start of the core making run, the shell corebox was preheated to and maintained at about 450°F. Fifteen cores were yielded for the core making portion of each run. An individual core took less than 2½ minutes to produce. Finished cores had the color of toast or

chestnut with an approximate wall thickness of 1/4 inch. An example of a sectioned finished shell core can be seen in Figure 2-3.

The start of the core making run commenced when sand was initially blown into the iron corebox. After the investment cycle the core was removed from the corebox, which signaled the end of the 2 minutes of core making time. The core was then immediately placed on a table located inside the enclosed hood for cooling and storage. This period was termed "core making/storage," because cooling cores were constantly present and potentially contributing

to emissions during core making. Cores were made continuously until the 15 cores were produced, which averaged just over 30 minutes. All cores were stored on the table under the hood immediately upon removal from the machine. After the 15<sup>th</sup> core was removed from the shell core machine and placed on the table along with the previously made cores, the separate core "storage" period began and continued until the predetermined 60 minute run time had elapsed. This separate storage period lasted just under 30 minutes.

# 2.1.3. Emissions Testing

Emissions testing included several methods for measuring both speciated and undifferentiated hydrocarbons and other chemical classes. The primary test method used for speciation was Method 18, one of the US Environmental Protection Agency's (EPA) reference methods for volatile organic compound (VOC) analysis from smokestacks and other industrial sources. Method 18 is an "umbrella" method, and is generally used to identify and/or measure as many compounds as possible in order to calculate actual VOC emissions from other measurements (e.g. EPA Method 25 or 25A). The method is a guideline and a system of quality assurance (QA) checks for VOC analysis rather than a rigorous, explicit manual for sampling or analysis.

As described in the method, sampling was conducted using a Volatile Organic Sampling Train (VOST). A sample gas stream was extracted from the source and then routed using the train through tubes containing adsorbents, which are the collection materials upon which the organic analytes are deposited. The selection of sampling media was based on the compounds expected to be evolved during the test. Adsorption tube samples were collected and analyzed for seventy-six (76) compounds using detailed collection and analysis procedures based on approved federal methods, including those of the EPA.

Two methods were employed to measure undifferentiated hydrocarbon emissions as Emission Indicators: TGOC as Propane, performed in accordance with EPA Method 25A, and non-methane hydrocarbons as determined from methane results obtained in a manner prescribed in EPA CTM-042.

Method 25A is an instrument based method in which the stack gas is introduced directly to a flame ionization detector (FID) without first separating the components. In Method 25A,

sampling is accomplished by extracting a gas stream from the stack effluent and transferring it via heated non-reactive tubing to the FID analyzer under very controlled temperature and pressure conditions. The FID measures the quantity of carbon containing molecules and is calibrated by a gas standard, which in this case is the three carbon alkane, propane  $(C_3H_8)$ . The FID will give a response relative to the calibration standard and results are expressed in terms of the gas used for calibration. Because the FID responds to all carbon containing compounds, methane  $(CH_4)$  and other exempt compounds are included in the total hydrocarbon results.

Methane was analyzed by a separate FID equipped with an oxidizing catalyst (methane cutter) that removes all non-methane hydrocarbons (NMHC). The calibration gas for this FID is methane (CH<sub>4</sub>). The two FIDs were run simultaneously, and collected data every second. NMHC results were then determined by subtracting the detected methane from the average total hydrocarbon value.

Continuous on-line monitoring of a subgroup of criteria pollutants and greenhouse gases including carbon dioxide (CO<sub>2</sub>), carbon monoxide (CO), and nitrogen oxides (NOx) was conducted according to US EPA Methods 3A, 10, and 7E, respectively.

Figure 2-4 shows the emission sampling equipment set-up for Test HO. The generated

Figure 2-4 Test Equipment for Emissions Sampling



emissions were transported through an insulated seven (7) inch stack located above the shell core machine. The flow rate of the emission capture air through the stack was nominally 220 scfm.

Heated sample probes inserted into the stack at relevant locations, determined by EPA Method 1, enabled collection of emissions directly from the stack. One probe provided gases for the sampling manifold to which was attached the VOST and a heated line for the methane analyzer. Another probe in the stack was used to continuously draw effluent samples and transport them through a second heated sample line to an emissions console that contains a battery of gas analyzers. This console, or emissions bench, consists of a total hydrocarbon analyzer for TGOC analysis, two infrared analyzers (for CO and CO<sub>2</sub>) and a chemiluminescence analyzer for NOx. All of the continuous monitors are located in Technikon's laboratory. Figure 2-5 shows details of the VOST and the emissions bench.

Figure 2-5 VOST and On-Line Analyzer

Volatile Organic Sampling Train

The heated sample lines were located approximately 120 feet away from the laboratory. A pressure drop of approximately 2 psi in the FID instruments resulted from the extended distance between the sampling location and the analytical instruments. Sample flows were decreased due to the extended distance, but were still within instrument ranges and sufficient for accurate sampling. Calibrations and calibration checks were performed throughout the test and indicated that the pressure drop could be compensated for, and did not significantly influence measurements.

Mass emission rates for all analytes were calculated using continuous monitoring or laboratory analytical results, measured source data and appropriate process data. Average emission results for individual runs are presented in Appendix B.

Speciated analyte emissions were calculated in addition to five "Emission Indicators:" TGOC as Propane, NMHC, Sum of Target Analytes, Sum of Target Hazardous Air Pollutants (HAPs), and the Sum of Target Polycyclic Organic Matter (POMs). Full descriptions of these indicators can be found in Section 3.0 of this report.

# 2.1.4. Process Parameter Measurements

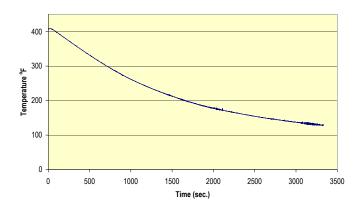
Table 2-2 lists the process parameters that were monitored during each test.

Table 2-2 Process Equipment and Methods

Process Parameters	Equipment and Method(s)
Core Weight	Mettler SB12001 Digital Scale (Gravimetric)
LOI%	Denver Instruments XE-100 Analytical Scale (AFS procedure 5100-00-S)
Core Box Temperature	Omega HH-23 Digital Thermometer

A thermocouple was placed inside of a cooling core during each of the six runs to track cooling rates, in addition to monitoring the core box temperature. Figure 2-6 gives the typical cooling decay rate from the maximum temperature reached by the core (approximately 405°F) to the end of the run, as exemplified by run HO003.

Figure 2-6 Cooling Rate of Shell
Core
H0003 Core Cooling



# 2.1.5. Sampling Methods

The specific sampling and analytical methods used for Test HO are based on federal reference methods shown in Table 2-3. The details of the specific testing procedures and their variance from the reference methods are included in the Technikon Standard Operating Procedures.

Table 2-3 Emission Sampling and Analytical Methods

Measurement Parameter	Test Method(s)
Port Location	US EPA Method 1
Number of Traverse Points	US EPA Method 1
Gas Velocity and Temperature	US EPA Method 2
Gas Density and Molecular Weight	US EPA Method 3a
Gas Moisture	US EPA Method 4 (Gravimetric)
Target Analytes including HAPs and POMs	US EPA Methods TO17, TO11; NIOSH
Target Analytes including HAPs and POWS	Methods S347; OSHA PV 2003
TGOC	US EPA Method 25A
CO	US EPA Method 10
CO <sub>2</sub>	US EPA Method 3A
NOx	US EPA Method 7E
SO <sub>2</sub>	OSHA ID 200
CH <sub>4</sub>	US EPA CTM 042

Some methods modified to meet specific CERP test objectives.

# 2.2. Data Reduction, Tabulation and Preliminary Report Preparation

Data calculations for determining emission concentrations resulting from the specific test plan outlined in Appendix A are based on process and emission parameters. The analytical results of the emissions sampling provide the mass of each analyte in the sample. The total mass of the analyte emitted is calculated by multiplying the mass of analyte in the sample by the ratio of total stack gas volume to sample volume. The total stack gas volume is calculated from the measured stack gas velocity and duct diameter and corrected to dry standard conditions using the measured stack pressures, temperatures, gas molecular weight and moisture content. The total mass of analyte is then divided by the weight of the casting poured or weight of binder to provide emissions data in pounds of analyte per ton core or pounds of analyte per pound of binder.

All emissions have been background subtracted. When sample measurements are made, the observed result includes the portion of the analyte in the sample, plus a response due to background contributions from either ambient or system blanks. The net analyte sample concentration is therefore the amount of the analyte, if any, found in the blank subtracted from the amount of analyte found in the sample. Background correcting the data allows

determination of emissions resulting only from the specific materials or process tested, and not those that may be present in either the ambient air of the research foundry, or from the manufacturing equipment used during the sampling period.

Subtracting those analytes found in quality control and method blanks (i.e. ambient background) results in emissions for the complete shell core making process, including all emissions from the machine, natural gas burners and any material emissions. Subtracting those analytes found in the system blank provides emissions for the materials used in making cores, including the resin. The system blank was obtained by operating the shell core machine for a simulated 60 minute run, but without core production. Any generated emissions were from operation of the shell core machine system only, including the 32 natural gas burners. Subtracting this system background from the analytical results obtained during testing eliminates all emissions given off by operation of the machine.

Individual speciated emission rates and reporting limit results for each analyte for all sampling runs are included in Appendix B of this report. Average emission results for the test are given in Section 3.0, Tables 3-1a and 3-1b.

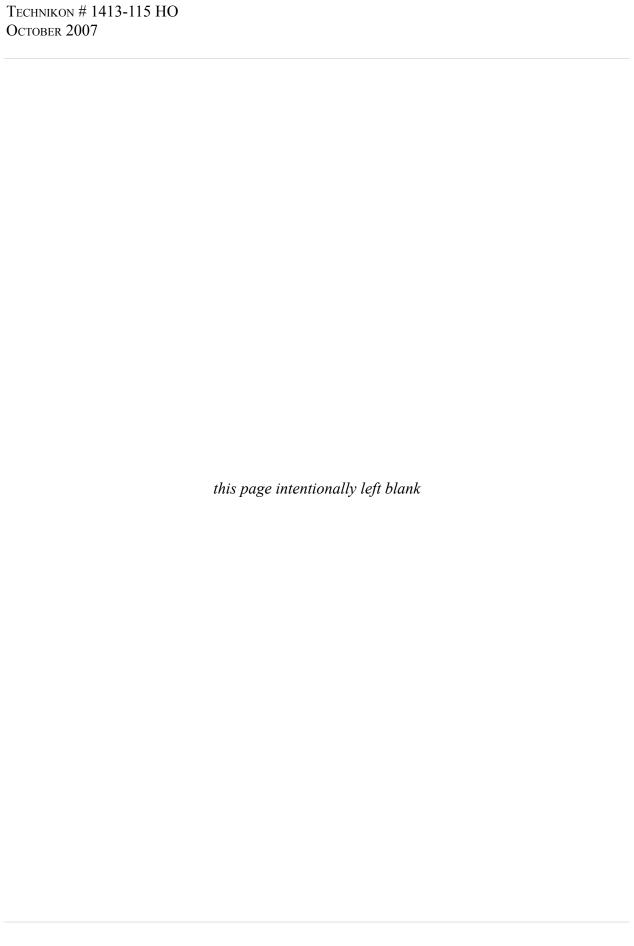
## 2.2.1. Report Preparation and Review

The Preliminary Draft Report is created and reviewed by Process Team and Emissions Team members to ensure its completeness, consistency with the test plan, and adherence to QA/QC procedures. Appropriate observations, conclusions and recommendations are added to the report to produce a Draft Report. The Draft Report is then reviewed by senior management and comments are incorporated into a draft Final Report prior to final signature approval and distribution.

# 2.3. Quality Assurance and Quality Control (QA/QC) Procedures

Detailed QA/QC and data validation procedures for the process parameters, stack measurements, and laboratory analytical procedures are included in the Technikon Emissions Testing and Analytical Testing Standard Operating Procedures. In order to ensure the timely review of critical quality control parameters, the following procedures are followed:

- Immediately following the individual sampling events performed for each test, specific process parameters are reviewed by the Process Engineer to ensure that the parameters are maintained within the prescribed control ranges. Where data are not within the prescribed ranges, the Manager of Process Engineering and the Vice President of Operations determine whether the individual test samples should be invalidated or flagged for further analysis following review of the laboratory data.
- The source (stack) and sampling parameters, analytical results and corresponding laboratory QA/QC data are reviewed by the Emissions Measurement Team to confirm the validity of the data. Senior management of Analytical Measurement Technologies reviews and approves the recommendation, if any, that individual sample data should be invalidated. Invalidated data are not used in subsequent calculations.



### 3.0 Test Results

Average results for the emissions measured during Test HO, and their comparison to the shell core making and storage baseline Test HN, are presented in Tables 3-1a and 3-1b. Concentrations are given in both lb/ton of cores and lb/lb of binder. All emission results presented in these tables were determined by subtracting the system background. The system background was obtained by operating the shell core machine without core production resulting in emissions from the materials used for core making and core storage only. Compounds found in the system background are reported in Table 3-2.

The individual chemical compounds targeted for collection and analyses from airborne emissions from the shell core process were chosen based on the chemistry of the binder under investigation, potential emissions from the operation of the shell core machine, and analytes historically targeted. These compounds, termed "target analytes" (TA), were chosen for their potential to emit based on chemical and foundry operational parameters.

Included in the tables are the "Sum of Target Analytes," the "Sum of Target HAPs," and the "Sum of Target POMs." These three analyte sums are part of the group termed "Emission Indicators." The emissions indicator called the "Sum of Target Analytes" is the sum of the individual speciated analytes that were targeted for collection and analysis, and detected at a level above the practical quantitation limit.

The target analyte sum includes targeted compounds that may also be defined as HAPs and POMs. By definition, HAPs are specific compounds listed in the Clean Air Act Amendments of 1990. The term POM defines a broad class of compounds based on chemical structure and boiling point. POMs as a class are a listed HAP. A subset of compounds from the current list of EPA HAPs was targeted for collection and analysis. These individual target HAPs (which may also be POMs by nature of their chemical properties) detected in the samples are summed together and defined as the "Sum of Target HAPs," while the "Sum of Target POMs" only sums those organic HAPs that are also defined as POMs.

Also included in this group and reported in the first section of the tables are "TGOC as

Propane" as determined by Method 25A, and NMHC. The second section of the tables includes average emission results for select HAPs and POMs. Additional individual target analytes are given in the third section of the tables. Average values for a subgroup of criteria and greenhouse gases including CO, CO<sub>2</sub>, CH<sub>4</sub>, SO<sub>2</sub>, and NO<sub>x</sub> are given in the fourth section of the tables. A few individual isomers are reported in the tables, and have not been combined or reported as a group. Isomers which have been targeted and analyzed may be summed as a group using the information located in these tables and in Appendix B.

Table 3-1a Test HO
Comparison Summary of
Selected Target Analytes to
Test HN, Average Results,
Ib/tn core

			Percent
	Reference		Change from
Analyte Name	Test HN	Test HO	Reference
Emission Indicators	169(1114	163(110	Reference
THC as Propane	5.34E-01	9.34E-01	75
Non-Methane Hydrocarbons	3.28E-01	2.96E-01	-10
Sum of Target Analytes	5.79E-01	1.46E-01	-75
Sum of Target HAPs	1.25E-01	4.06E-02	-67
Sum of Target POMs	ND	4.69E-05	NA
Selected Target HAPs and PO		11002 00	100
Phenol	1.12E-01	3.23E-02	-71
Formaldehyde	1.11E-02	7.18E-03	-35
Toluene	4.46E-04	3.27E-05	-93
Propionaldehyde (Propanal)	4.30E-04	1.79E-04	-58
Acetaldehyde	3.62E-04	2.62E-04	-28
Acrolein	1.17E-04	1.54E-04	32
Benzene	≤PQL	2.95E-05	NA
Cresol, o-	≤PQL	3.16E-05	NA
Hexane	≤PQL	6.38E-05	NA
Naphthalene	≤PQL	4.69E-05	NA
Styrene	≤PQL	3.01E-04	NA
Additional Selected Target An	alytes		
Ammonia	4.53E-01	1.03E-01	-77
Crotonaldehyde	4.81E-04	8.70E-04	81
2-Butanone (MEK)	3.35E-04	3.02E-04	-10
Benzaldehyde	2.47E-04	≤PQL	NA
Butyraldehyde/Methacrolein	2.46E-04	2.94E-04	20
Pentanal (Valeraldehyde)	1.15E-04	≤PQL	NA
Butylbenzene, sec-	≤PQL	2.99E-05	NA
Diethylbenzene, 1,3-	≤PQL	3.02E-05	NA
Diethylbenzene, 1,4-	≤PQL	3.74E-05	NA
Dimethylphenol, 3,4-	≤PQL	8.01E-05	NA
Heptane	≤PQL	4.21E-05	NA
Tetradecane	≤PQL	2.87E-05	NA
Trimethylbenzene, 1,2,4-	≤PQL	4.60E-05	NA
Criteria Pollutants and Greenhouse Gases			
Carbon Dioxide	1.59E+01	1.49E+01	-6
Methane	2.06E-01	6.38E-01	210
Carbon Monoxide	1.26E-01	1.15E-01	-9
Nitrogen Oxides	≤PQL	≤PQL	NA
Sulfur Dioxide	≤PQL	≤PQL	NA

ND=Not Detected

NA= Not Applicable

I=Invalidated Data

 $\leq$ PQL=Less than or equal to the Practical Quantitation Limit

Table 3-1b Test HO
Comparison Summary of
Selected Target Analytes to
Test HN, Average Results, lb/
Ib binder

	<u> </u>		Percent	
			Change	
	Reference		from	
Analyte Name	Test HN	Test HO	Reference	
Emission Indicators				
THC as Propane	9.16E-03	1.60E-02	75	
Non-Methane Hydrocarbons	5.63E-03	5.08E-03	-10	
Sum of Target Analytes	1.00E-02	2.53E-03	-75	
Sum of Target HAPs	2.17E-03	7.07E-04	-67	
Sum of Target POMs	ND	8.15E-07	NA	
Selected Target HAPs and PO		0.102 07	1471	
Phenol	1.95E-03	5.62E-04	-71	
Formaldehyde	1.93E-04	1.25E-04	-35	
Toluene	7.79E-06	5.64E-07	-93	
Propionaldehyde (Propanal)	7.48E-06	3.17E-06	-58	
Acetaldehyde	6.31E-06	4.58E-06	-27	
Acrolein	2.03E-06	2.72E-06	34	
Naphthalene	≤PQL	8.15E-07	NA	
Benzene	≤PQL	5.13E-07	NA	
Cresol, o-	≤PQL	5.42E-07	NA	
Hexane	≤PQL	1.11E-06	NA	
Styrene	≤PQL	5.24E-06	NA	
Additional Selected Target Analytes				
Ammonia	7.85E-03	1.80E-03	-77	
Crotonaldehyde	8.37E-06	1.53E-05	83	
2-Butanone (MEK)	5.82E-06	5.23E-06	-10	
Benzaldehyde	4.30E-06	≤PQL	NA	
Butyraldehyde/Methacrolein	4.27E-06	5.24E-06	23	
Pentanal (Valeraldehyde)	1.98E-06	≤PQL	NA	
Tetradecane	≤PQL	4.95E-07	NA	
Diethylbenzene, 1,3-	≤PQL	5.19E-07	NA	
Heptane	≤PQL	7.33E-07	NA	
Trimethylbenzene, 1,2,4-	≤PQL	8.00E-07	NA	
Cymene, p-	≤PQL	4.84E-07	NA NA	
Butylbenzene, sec-	≤PQL	5.13E-07	NA NA	
Diethylbenzene, 1,4-	≤PQL ≤PQL	6.42E-07	NA NA	
Dimethylphenol, 3,4- Criteria Pollutants and Greenl		1.37E-06	NA	
Carbon Dioxide	2.73E-01	2.56E-01	-6	
Methane	3.54E-03	1.10E-02	210	
Carbon Monoxide	2.16E-03	1.10L-02 1.97E-03	-9	
Nitrogen Oxides	≤PQL	≤PQL	NA	
Sulfur Dioxide	≤PQL	≤PQL	NA	
- Canar Dioxido	<-	&-	. */ `	

ND=Not Detected

NA= Not Applicable

I=Invalidated Data

≤PQL=Less than or equal to the Practical Quantitation Limit

Compounds found in the system background are reported in Table 3-2. These values were subtracted from the analytical results for each run prior to calculating emission factors in units of lb/lb binder and lb/ton cores.

Table 3-2 System Background Emissions for Test HO

Analyte	Value	Unit
TGOC as Propane	109.4	ppm
CO	14.6	ppm
CO <sub>2</sub>	1381.5	ppm
NO <sub>x</sub>	0.5	ppm
Acenaphthalene	49.3	ng
Anthracene	113.0	ng
Benzene	33.4	ng
Butylbenzene, n-	28.2	ng
Butylbenzene, tert-	201.0	ng
Cyclohexane	47.5	ng
Decane	24.2	ng
Diisopropylbenzene, 1,3-	85.2	ng
Dimethylnaphthalene, 1,3-	25.5	ng
Ethylbenzene	34.0	ng
Heptane	53.6	ng
Hexane	138.0	ng
Methylnaphthalene, 1-	215.0	ng
Methylnaphthalene, 2-	25.4	ng
Octane	39.1	ng
Phenol	249.0	ng
Styrene	34.4	ng
Toluene	59.8	ng
Trimethylbenzene, 1,2,3-	30.6	ng
Xylene, mp-	90.2	ng
Xylene, o-	27.7	ng
Acetaldehyde	389.0	μg
Formaldehyde	6580.0	μg

Emissions data that have been determined to be below the practical quantitation limit (PQL) after data validation and verification are substituted with the numerical value used for the PQL, rather than with the value of zero. This practice results in a conservative value for the average test emissions. If an analyte has calculated concentrations above the PQL for some runs, but values for other runs fall below the PQL, the PQL value is included when calculating analyte averages and sums. However, if an analyte has a concentration that is below the PQL for <u>all</u> runs in a test, the test average is indicated by ≤PQL (less than or equal to the PQL) in the Tables and Figures of this report, and no runs are included in any summations or averages. Omitting these less-than-reporting-limit analytes in calculations ensures that only those targeted compounds which contribute to emissions are included in emission sums.

Examination of measured process parameters indicated that Test HO was run within acceptable ranges and limits. The primary causes and secondary influences on emissions were controlled for each individual run, so that for shell core making and storage, resultant air emissions reflect only the difference in the process and materials being tested.

### 3.1. Discussion of Results

Figures 3-1a to 3-4a (lb/ton of core) and Figures 3-1b to 3-4b (lb/lb of binder) graphically present the emissions data from Tables 3-1a and 3-1b for Test HO compared to Test HN for the five emissions indicators, selected individual HAP, other target analyte, and selected criteria pollutant and greenhouse gases.

Figure 3-1a Comparison of Emissions Indicators of Test HO to Reference Test HN, Average Results, lb/tn core

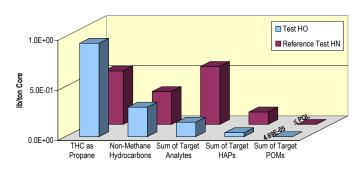


Figure 3-2a Comparison of Selected HAP and POM Emissions of Test HO to Reference Test HN, Average Results, lb/tn core

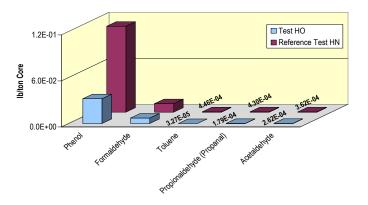


Figure 3-3a Comparison of Selected Target Analyte Emissions of Test HO to Reference Test HN, Average Results, lb/tn core

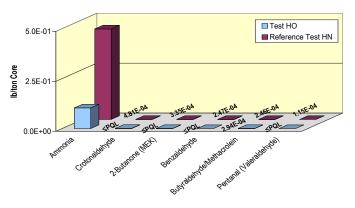


Figure 3-4a Comparison of Criteria Pollutants and Greenhouse Gases of Test HO to Reference Test HN, Average Results, Ib/tn core

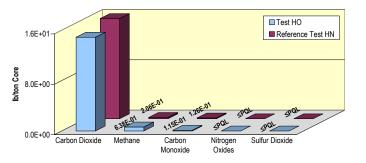


Figure 3-1b Comparison of Emissions Indicators of Test HO to Reference Test HN, Average Results, Ib/lb binder

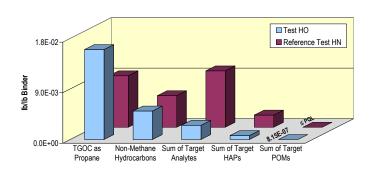


Figure 3-2b Comparison of Selected HAP and POM Emissions of Test HO to Reference Test HN, Average Results, lb/lb binder

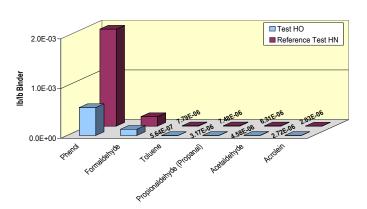


Figure 3-3b Comparison of Selected Target Analyte Emissions of Test HO to Reference Test HN, Average Results, Ib/lb binder

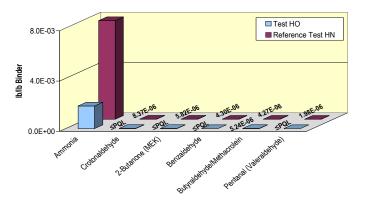
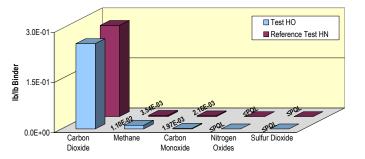


Figure 3-4b Comparison of Criteria Pollutants and Greenhouse Gases of Test HO to Reference Test HN, Average Results, lb/lb binder



Two target analytes comprised over 97% of emissions for the system background adjusted data. Ammonia had the highest contribution at 78% followed by phenol at 19%. Formaldehyde was the third highest emitter, but contributed only 2% to emissions. Phenol and formaldehyde are HAPs, and although ammonia is not an EPA regulated HAP, it is a reportable chemical as part of the EPA's EPCRA Section 313 chemical list. The few compounds which supplied the remainder of measurable emissions were all less than 0.1%. Of the 76 target analytes, 22 were responsible for emissions, including 11 HAPs. Targeted HAPs contributed approximately 22% to total measured emissions.

Five target analyte emissions were lower and two were higher when comparing the two binder systems used in Test HO and Test HN. The largest emissions decrease of 77% was found for ammonia, although its contribution (78%) to the total measured targeted emissions was exactly the same as in Test HN. The largest increase in emissions of 81% was for crotonaldehyde.

The target analyte sum would theoretically closely match the results for total hydrocarbons obtained by Method 25A, excluding exempt compounds such as methane, and including compounds such as formaldehyde and ammonia, which are less responsive in the FID. For the results reported here, the Sum of Target Analytes is only 17% of the measured TGOC as Propane results. Calculating TGOC as Propane value adjusted by the above named compounds gives a value that is only 12% lower than the measured TGOC as Propane value.

A GC/MS analysis of the stack emissions revealed several additional compounds for both Test HO and Test HN. Some of these compounds were of higher concentration than the predicted compounds, and several are on either the US EPA's HAP or SARA 313 chemical lists. A probability based matching algorithm was used to match the mass spectra of the unknown compounds to a standard GC/MS library (NIST Mass Spectral Database, 1998). Tentative identification included several silicone compounds (likely from the release agent used during core making) and numerous nitrogen containing compounds, including nitro phenols. A list of these compounds is given in Table 3-3.

Table 3-3 Additional Identified Compounds from GC/MS Analysis

Compound		
3-Penten-2-Ol		
4H-3,1-Benzoxazin-2(1H)-One		
Acetonitrile, Dimethylamino		
Benzaldehyde, 2-Hydroxy		
Benzeneacetonitrile, 4-Hydroxy-		
Benzoic Acid, 2-[(Trimethylsilyl)O		
Cyclotetrasiloxane, Octamethyl-		
Cyclotrisiloxane, Hexamethyl		
Ethylenimine #		
Formamide, N,N-Dimethyl- <sup>+</sup>		
O-(N-Methylformimidoyl)Phenol		
P-Benzoquinone <sup>#</sup>		
Pentane		
Pentane, 2-Methyl- 2-Meth		
Phenol, 2-Nitro- <sup>+</sup>		
Phenol, 4-Methyl-2-Nitro-		
Phenol, 4-Methyl-2-Nitro-		
Silane, 5,5-Dimethyl-4-Methylene		
Silicone Polymer		
Silicone Polymer		
Spiro[2.4]Hept-5-Ene, 5-Trimethyls		

<sup>#</sup> HAP Chemical

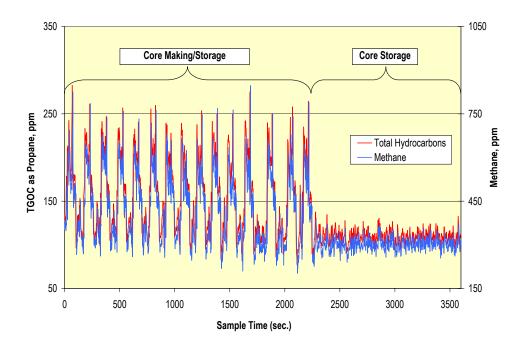
Material emissions for core making/storage and storage were attained by subtracting a background sample obtained by operating the shell core machine without core production. The system background contained moderate concentrations of numerous hydrocarbons including benzene, toluene, phenol, and xylenes. The measured criteria pollutants and greenhouse gases concentration were also high in the system blank. Comparing system blank concentrations to sample concentrations shows that a majority of the differentiated hydrocarbons, the total hydrocarbons as measured by TGOC as Propane, and the CH<sub>4</sub> originated from the shell core machine, including the natural gas burners. It is also apparent that

<sup>\*</sup> SARA 313 Chemical

most of the total hydrocarbon emissions as measured by TGOC as propane is comprised of CH<sub>4</sub>. The emissions for total hydrocarbons and methane parallel each other exactly over the entire core making/storage and storage run, as shown by example for run HO003 in Figure 3-5.

Figure 3-5 Comparison of Methane and TGOC Results for Shell Core Making

Run H0003



For the continuous monitors, data were acquired every second, and the average value calculated over the entire sampling period is reported as the run average. For both Test HO and Test HN, this average value includes data acquired during the actual core making as well as the short interval between cores (after a finished core was removed from the machine until another was made), and the final 30 minute core storage period (where the cooling cores were stored under the hood after the core making/storage period). The separate storage and cooling period that began after the removal of the fifteenth core shows low emissions. The average for each run is therefore an underestimation of emissions. An example of the cyclical nature of the core making and storage is shown in Figure 3-6 for  $\mathrm{CO}_2$ .

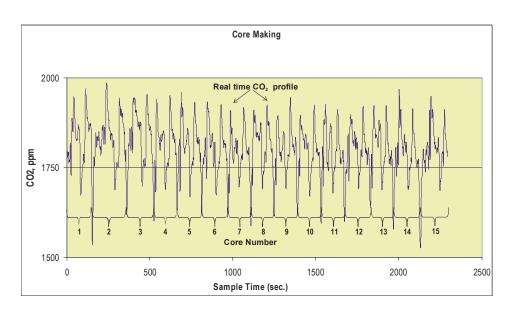


Figure 3-6 Cyclical Pattern of Emissions from Shell Core Making

Stack samples were also continuously drawn through the media on the VOST during the entire cyclical core making/storage and separate storage periods. The sampling media is minimally affected by those intervals between cores and the separate storage period when there were few to no emissions. This is due to the sample being one single sample that is consolidated over the entire run, rather than a calculated average of numerous samples.

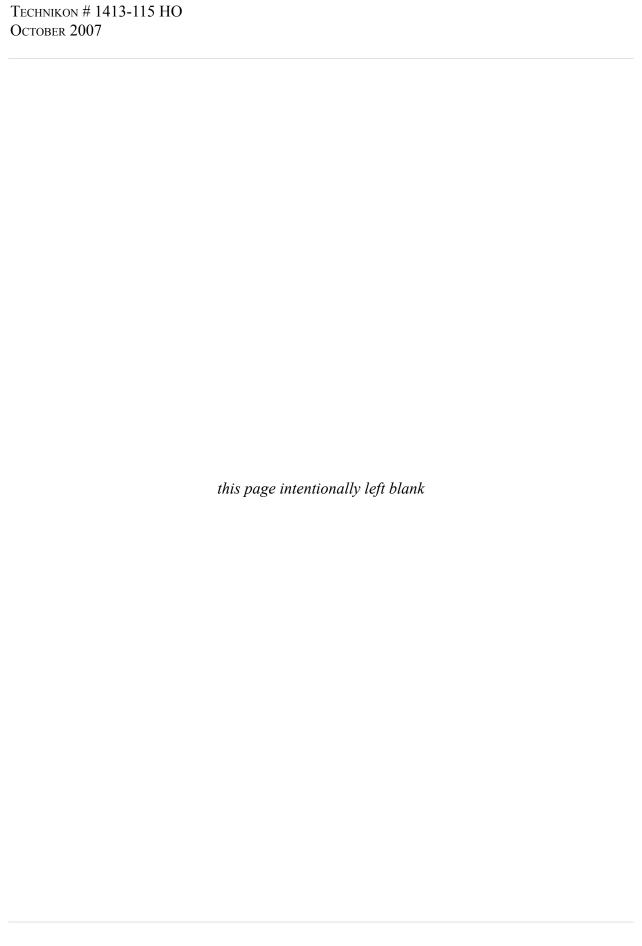
The cooling curve for the cores was anticipated to mimic the decrease in emissions on core storage. However, in comparing Figure 2-6 with the storage period shown in Figure 3-5, emissions decreased much faster than the cooling curve. Immediately upon removal from the machine, emissions dropped almost to background levels.

Average core parameters for Test HO are summarized in Table 3-4, and detailed information regarding testing, sampling, data collection and results for each sampling event are contained in the four appendices. Appendix A contains test plans, instructions and the sampling plans for Test HO. Appendix B contains detailed emissions data and average results for all targeted analytes. Target analyte practical quantitation limits expressed in both lb/lb binder and lb/ton metal are also shown in Appendix B. Appendix C contains detailed process data. Appendix D contains continuous monitor charts. The charts are pre-

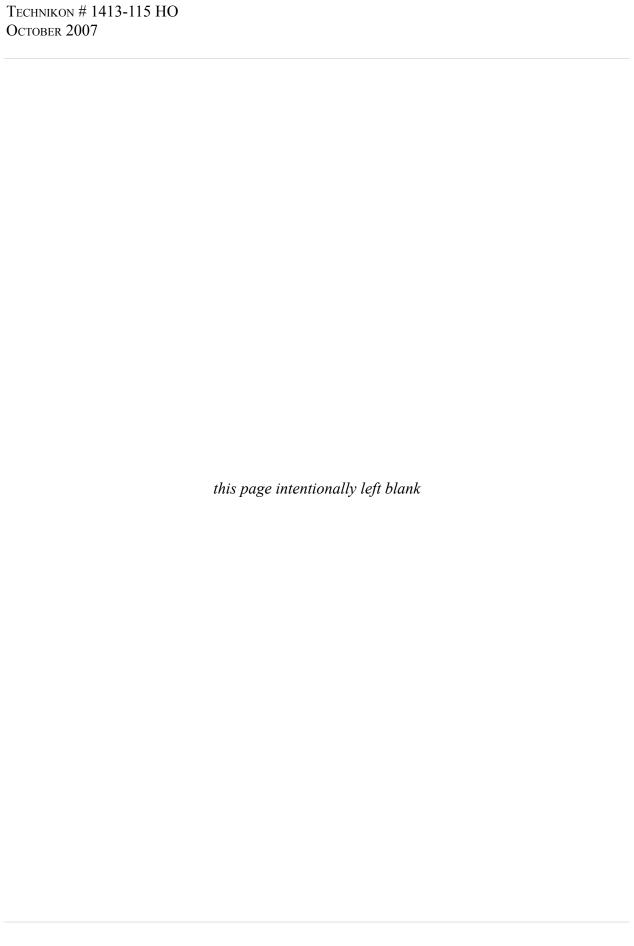
sented to show TGOC, carbon monoxide, carbon dioxide, methane, and oxides of nitrogen time-dependent emissions profiles for each individual emissions run. Charts have not been background corrected. Appendix E contains acronyms and abbreviations.

Table 3-4 Summary of Core Making Process Parameters

	Test HN	Test HO
Test Dates	4/23/07-4/25/07	4/25/07-4/26/07
Average cooled core weight, g	2515.6	2617.9
Average core box temperature	456	448
% stated core binder (BOS)	3	3
% calculated actual binder	2.9	2.9
Number of cores made in run	15	15
Total core weight in from run, lbs.	83.2	86.6
Total binder weight in run, lbs.	2.4	2.5
Core LOI, %	3.21	2.86



## APPENDIX A TEST AND SAMPLE PLANS AND PROCESS INSTRUCTIONS



## **Technikon Test Plan** page 1 of 2 Fill-in and check all that apply 1413 TASK NUMBER 122 DOUBLE ALPHA HN CONTRACT NUMBER: On Site at Research Foundry Off Site at • SITE: • DATE RANGE: From Apr. 23, 2007 to APR. 25, 2007 Emissions Testing Mechanical Properties Casting Quality • TEST TYPE: Baseline Comparison to \_\_\_\_ Other \_\_\_ Core Mold • PROCESSES: Pouring Cooling Shakeout Mixing Making Storage ☐ Iron ☐ Aluminum ☐ Steel ☐ Other \_\_\_\_ Pour Temp: \_\_\_\_°F NA • METAL: Step Star Irregular Gear Other • PATTERN: • MOLD: Number Molds \_\_\_\_\_ Number Cavities NA Flask Dimensions \_\_\_\_\_ Storage Temp: \_\_\_\_\_°F Cold box Warm box Hot box No-bake Shell Oil • CORE TYPE: Phenolic Urethane Furan Inorganic (inc. Sodium Silicate) Epoxy-Acrylic Alkaline Phenolic Ester Coated Core Wash Naphthalene Depleted CO2 Cured SO2 Cured Acid Cured TEA Cured Methyl Formate Cured Hot Air Cured Other \_\_\_\_ Product Name(s) Technisand XC30 • CORE COATING: None All Runs Conditioning Runs Only Test Runs Only Baumé Application Method \_\_\_\_\_ Drying Method \_\_\_\_\_ Product Name(s) \_\_\_\_ Greensand Seacoal No-Bake Other Shell • CORE SAND: Additives \_\_\_\_ Product Name(s) Technisand XC 30 • CORE BINDER: Type Phenol-Formaldehyde Concentration 3% (BOS) Ratio $(\frac{P1}{P2})$ Product Name(s) Novalac Type NA Concentration BOS BOR • ACTIVATOR: Product Name(s) \_\_\_\_

## **Technikon Test Plan**

page 2 of 2

		Fill-in and che	eck all that apply		
CONTRACT NUMBER:	1413	TASK NUMBER	122	DOUBLE ALPHA	HN
RELEASE AGENT:	Type Concentropy Product Name(s) Ashle		•		
Runs:	Number for Conditioni	ng Duration	on Description _		ng per core for a total of
WASTE MATERIAL TO BE SAMPLED:	All None For: Contaminates list	<u> </u>	_	og bones	
TEST OBJECTIVES:	To measure emissions sand.	s from shell core n	naking of Fairmount mi	inerals Technisand	XC30 3% BOS shell
RESULTS TO BE REPORTED:			ect Target Analytes and sper ton of cores ma		orted in pounds of
ADDITIONAL COMMENTS:			ne to perform this test. o operate the machine.		otal enclosure of the

This test plan is routed to and/or reviewed by the following:

Senior Process Engineer Production Engineer Measurement Technology Manager V.P. Operations Steering Committee Emissions Team Chair

## **Technikon Test Plan** page 1 of 2 Fill-in and check all that apply 1413 TASK NUMBER 115 DOUBLE ALPHA HO CONTRACT NUMBER: On Site at Research Foundry Off Site at \_\_\_\_\_ • SITE: • DATE RANGE: From Apr 25, 2007 to Apr 26, 2007 Emissions Testing Mechanical Properties Casting Quality • TEST TYPE: Baseline X Comparison to HN Other Core Mold • PROCESSES: Pouring Cooling Shakeout Mixing Making Storage ☐ Iron ☐ Aluminum ☐ Steel ☐ Other \_\_\_\_ Pour Temp: \_\_\_\_°F NA • METAL: Step Star Irregular Gear Other • PATTERN: • MOLD: Number Molds \_\_\_\_\_ Number Cavities \_\_\_\_\_ NA Flask Dimensions \_\_\_\_\_ Storage Temp: \_\_\_\_\_°F Cold box Warm box Hot box No-bake Shell Oil • CORE TYPE: Phenolic Urethane Furan Inorganic (inc. Sodium Silicate) Epoxy-Acrylic Alkaline Phenolic Ester Coated Core Wash Naphthalene Depleted CO<sub>2</sub> Cured SO<sub>2</sub> Cured Acid Cured TEA Cured Methyl Formate Cured Hot Air Cured Other \_\_\_\_ Product Name(s) Technisand 630BN Gold None All Runs Conditioning Runs Only Test Runs Only Baumé • CORE COATING: Application Method \_\_\_\_\_ Drying Method \_\_\_\_\_ Product Name(s) Greensand Seacoal No-Bake Other Shell + CORE SAND: Additives Product Name(s) Technisand 630BN Gold Type Phenol-Formaldehyde Concentration 3% (BOS) Ratio ( $\frac{P1}{P2}$ )\_\_\_\_\_ • CORE BINDER: Product Name(s) Novalac Type NA Concentration \_\_\_\_ BOS BOR • ACTIVATOR: Product Name(s) \_\_\_\_

Measurement Technology Manager

Steering Committee Emissions Team Chair

V.P. Operations

## **Technikon Test Plan** page 2 of 2 Fill-in and check all that apply 1413 TASK NUMBER 115 НΟ CONTRACT NUMBER: DOUBLE ALPHA Type \_\_\_\_ Concentration \_\_\_\_ (BOS) **◆ RELEASE AGENT:** Product Name(s) Ashland Zip-Slip 109 W • Runs: Number for Conditioning Duration Description Number for Sampling 6 Duration 15 Description 3 minutes making per core for a total of 1 hour sampling time per test All None Core Butts Loose Sand Dog bones **♦ WASTE MATERIAL** TO BE SAMPLED: For: Contaminates listed in EPA CRF section 40 part 261.24 To measure emissions from shell core making of Fairmount minerals Technisand 630BN Gold 3% BOS • TEST OBJECTIVES: shell sand. • RESULTS TO BE CO, CO2, NOx, CH4, NH3, SO2 and Select Target Analytes and POMs to be reported in pounds of emissions per pound of binder and pounds per ton of cores made. REPORTED: ADDITIONAL Technikon will use the shell core machine to perform this test. There will not be total enclosure of the COMMENTS: machine in order to allow the operator to operate the machine.. This test plan reviewed by: Senior Process Engineer **Production Engineer**

## **RESEARCH FOUNDRY HN - SERIES SAMPLE PLAN**

RESEARCH FOUNDRY HIN - SERIES SAMPLE PLAN													
Method	Sample #	Data	Sample	Duplicate	Blank	Breakthrough	Spike	Spike Duplicate	Flow (ml/min)		Comments: One hour of core making accompanied by an additional half hour of core storage for a total period of 90 minutes		
4/25/2007													
THC, CH4, CO, CO <sub>2</sub> & NOx	HN010	Х									TOTAL		
TO-17	HN01001		1						60	1	Carbopak charcoal		
	Excess								60	2	BLOCKED		
	Excess								60	3	BLOCKED		
Acetophenone	HN01002		1						200	4	15/30 mg Tenax (SKC 226-35-03)		
	Excess								200	5	BLOCKED		
NIOSH S347	HN01003		1						1000	6	Acid Silica Gel (SKC 226-10-06)		
NIOSH 6016	HN01004		1						1000	7	Acid Silica Gel (SKC 226-10-06)		
OSHA ID200	HN01005		1						1000	8	100/50 mg Carbon Bead (SKC 226-80)		
	Excess								1000	9	BLOCKED		
TO11	HN01006		1						1000	10	DNPH Silica Gel (SKC 226-119)		
	Excess								1000	11	BLOCKED		
	Moisture		1						500	12	TOTAL		
	Excess								5000	13	Excess		

## RESEARCH FOUNDRY HN - SERIES SAMPLE PLAN

Method	Sample #	Data	Sample	Duplicate	Blank	Breakthrough	Spike	Spike Duplicate	Flow (ml/min)		Comments: System blank with core blower burners on
THC, CH4, CO, CO <sub>2</sub> & NOx	HN011	Х									TOTAL
TO-17	HN01101		1						60	1	Carbopak charcoal
	Excess								60	2	BLOCKED
	Excess								60	3	BLOCKED
Acetophenone	HN01102		1						200	4	15/30 mg Tenax (SKC 226-35-03)
	Excess								200	5	BLOCKED
NIOSH S347	HN01103		1						1000	6	Acid Silica Gel (SKC 226-10-06)
NIOSH 6016	HN01104								1000	7	Acid Silica Gel (SKC 226-10-06)
OSHA ID200	HN01105		1						1000	8	100/50 mg Carbon Bead (SKC 226-80)
	Excess								1000	9	BLOCKED
TO11	HN01106		1						1000	10	DNPH Silica Gel (SKC 226-119)
	Excess								1000	11	BLOCKED
	Moisture		1						500	12	TOTAL
	Excess								5000	13	Excess

## **RESEARCH FOUNDRY HN - SERIES SAMPLE PLAN**

	Sample #		ple	Duplicate	k	Breakthrough	9	e Duplicate	Flow (ml/min)	Channel	
Method	sam	Data	Sample	ᆵ	Blank	3rea	Spike	Spike	<u>8</u>	Train (	Comments:
4/23/2007	0,		0,				0,	0,			Comments:
THC, CH4, CO, CO <sub>2</sub> & NOx	HN001	Х									TOTAL
TO-17	HN00101		1						60	1	Carbopak charcoal
TO-17	HN00102				1				0		Carbopak charcoal
	Excess								60	2	BLOCKED
	Excess								60	3	BLOCKED
Acetophenone	HN00103		1						200	4	15/30 mg Tenax (SKC 226-35-03)
Acetophenone	HN00104				1				0		15/30 mg Tenax (SKC 226-35-03)
	Excess								200	5	BLOCKED
NIOSH S347			1						1000	6	Acid Silica Gel (SKC 226-10-06)
NIOSH S347	HN00106				1				0		Acid Silica Gel (SKC 226-10-06)
NIOSH 6016	HN00107		1						1000	7	Acid Silica Gel (SKC 226-10-06)
NIOSH 6016	HN00108				1				0		Acid Silica Gel (SKC 226-10-06)
OSHA ID200	HN00109		1						1000	8	100/50 mg Carbon Bead (SKC 226-80)
OSHA ID200	HN00110				1				0		100/50 mg Carbon Bead (SKC 226-80)
TO11			1						1000	9	DNPH Silica Gel (SKC 226-119)
TO11				1					1000	10	DNPH Silica Gel (SKC 226-119)
TO11	HN00113				1				0		DNPH Silica Gel (SKC 226-119)
	Excess								1000	11	BLOCKED
	Moisture		1						500	12	TOTAL
	Excess								5000	13	Excess

#### RESEARCH FOUNDRY HN - SERIES SAMPLE PLAN

RESEARCH FOUNDRY HN - SERIES SAMPLE PLAN													
Method	Sample #	Data	Sample	Duplicate	Blank	Breakthrough	Spike	Spike Duplicate	Flow (ml/min)	Train Channel	Comments:		
4/23/2007													
THC, CH4, CO, CO <sub>2</sub> & NOx	HN002	Х									TOTAL		
TO-17	HN00201		1						60	1	Carbopak charcoal		
TO-17	HN00202			1					60	2	Carbopak charcoal		
	Excess								60	3	BLOCKED		
Acetophenone	HN00203		1						200	4	15/30 mg Tenax (SKC 226-35-03)		
Acetophenone	HN00204			1					200	5	15/30 mg Tenax (SKC 226-35-03)		
NIOSH S347	HN00205		1						1000	6	Acid Silica Gel (SKC 226-10-06)		
NIOSH S347	HN00206			1					1000	7	Acid Silica Gel (SKC 226-10-06)		
NIOSH 6016	HN00207		1						1000	8	Acid Silica Gel (SKC 226-10-06)		
OSHA ID200	HN00208		1						1000	9	100/50 mg Carbon Bead (SKC 226-80)		
	Excess								1000	10	BLOCKED		
TO11	HN00209		1						1000	11	DNPH Silica Gel (SKC 226-119)		
	Moisture		1						500	12	TOTAL		
	Excess								5000	13	Excess		

## **RESEARCH FOUNDRY HN - SERIES SAMPLE PLAN**

Method	Sample #	Data	Sample	Duplicate	Blank	Breakthrough	Spike	Spike Duplicate	Flow (ml/min)	Train Channel	Comments:
4/24/2007											
THC, CH4, CO, CO <sub>2</sub> & NOx	HN003	Х									TOTAL
TO-17	HN00301		1						60	1	Carbopak charcoal
TO-17 MS	HN00302		1						60	2	Carbopak charcoal
TO-17 MS	HN00303			1					60	3	Carbopak charcoal
	Excess								200	4	BLOCKED
Acetophenone	HN00304		1						200	5	15/30 mg Tenax (SKC 226-35-03)
NIOSH S347	HN00305		1						1000	6	Acid Silica Gel (SKC 226-10-06)
NIOSH 6016	HN00306		1						1000	7	Acid Silica Gel (SKC 226-10-06)
NIOSH 6016	HN00307			1					1000	8	Acid Silica Gel (SKC 226-10-06)
OSHA ID200	HN00308		1						1000	9	100/50 mg Carbon Bead (SKC 226-80)
OSHA ID200	HN00309			1					1000	10	100/50 mg Carbon Bead (SKC 226-80)
TO11	HN00310		1						1000	11	DNPH Silica Gel (SKC 226-119)
	Moisture		1						500	12	TOTAL
	Excess								5000	13	Excess

DECEMBALL FOLINDS	DV LINE OFFICE	OAMBLE BLAN
RESEARCH FOUND	RY HN - SERIES	SAMPLE PLAN

RESEARCH FOORBILL THE SERVED GAME EET EAR													
Method	Sample #	Data	Sample	Duplicate	Blank	Breakthrough	Spike	Spike Duplicate	Flow (ml/min)	Train Channel	Comments:		
4/24/2007													
THC, CH4, CO, CO <sub>2</sub> & NOx	HN004	Х									TOTAL		
TO-17	HN00401		1						60	1	Carbopak charcoal		
TO-17	HN00402					1			60	1	Carbopak charcoal		
	Excess								60	2	BLOCKED		
	Excess								60	3	BLCOKED		
Acetophenone	HN00403		1						200	4	15/30 mg Tenax (SKC 226-35-03)		
	Excess								200	5	BLOCKED		
NIOSH S347	HN00404		1						1000	6	Acid Silica Gel (SKC 226-10-06)		
	Excess								1000	7	BLOCKED		
OSHA ID200	HN00405		1						1000	8	100/50 mg Carbon Bead (SKC 226-80)		
	Excess								1000	9	BLOCKED		
TO11	HN00406		1						1000	10	DNPH Silica Gel (SKC 226-119)		
	Excess								1000	11	BLOCKED		
	Moisture		1						500	12	TOTAL		
	Excess								5000	13	Excess		

## **RESEARCH FOUNDRY HN - SERIES SAMPLE PLAN**

KECEAROTT CORD											
Method	Sample#	Data	Sample	Duplicate	Blank	Breakthrough	Spike	Spike Duplicate	Flow (ml/min)	Train Channel	Comments:
4/24/2007											
THC, CH4, CO, CO <sub>2</sub> & NOx	HN005	Х									TOTAL
TO-17	HN00501		1						60	1	Carbopak charcoal
	Excess								60	2	BLOCKED
	Excess								60	3	BLOCKED
Acetophenone	HN00502		1						200	4	15/30 mg Tenax (SKC 226-35-03)
	Excess								200	5	BLOCKED
NIOSH S347	HN00503		1						1000	6	Acid Silica Gel (SKC 226-10-06)
	Excess								1000	7	BLOCKED
OSHA ID200	HN00504		1						1000	8	100/50 mg Carbon Bead (SKC 226-80)
	Excess								1000	9	BLOCKED
TO11	HN00505		1						1000	10	DNPH Silica Gel (SKC 226-119)
	Excess								1000	11	BLOCKED
	Moisture		1						500	12	TOTAL
	Excess								5000	13	Excess

#### **RESEARCH FOUNDRY HN - SERIES SAMPLE PLAN**

RESEARCH FOUNDRY HIN - SERIES SAMPLE PLAN													
Method	Sample #	Data	Sample	Duplicate	Blank	Breakthrough	Spike	Spike Duplicate	Flow (ml/min)		Comments:		
4/25/2007													
THC, CH4, CO, CO <sub>2</sub> & NOx	HN006	Х									TOTAL		
TO-17	HN00601		1						60	1	Carbopak charcoal		
	Excess								60	2	BLOCKED		
	Excess								60	3	BLOCKED		
Acetophenone	HN00602		1						200	4	15/30 mg Tenax (SKC 226-35-03)		
	Excess								200	5	BLOCKED		
NIOSH S347	HN00603		1						1000	6	Acid Silica Gel (SKC 226-10-06)		
	Excess								1000	7	BLOCKED		
OSHA ID200	HN00604		1						1000	8	100/50 mg Carbon Bead (SKC 226-80)		
	Excess								1000	9	BLOCKED		
TO11	HN00605		1						1000	10	DNPH Silica Gel (SKC 226-119)		
	Excess								1000	11	BLOCKED		
	Moisture		1						500	12	TOTAL		
	Excess								5000	13	Excess		

## **RESEARCH FOUNDRY HO - SERIES SAMPLE PLAN**

							—				
Method	Sample #	Data	Sample	Duplicate	Blank	Breakthrough	Spike	Spike Duplicate	Flow (ml/min)		Comments: One hour of core making accompanied by an additional half hour of core storage for a total period of 90 minutes. TEST NOT DONE.
THC, CH4, CO, CO <sub>2</sub> & NOx	HO010	Х									TOTAL
TO-17	HO01001		1						200	1	Carbopak charcoal
	Excess								200	2	BLOCKED
	Excess								200	3	BLOCKED
	Excess								400	4	BLOCKED
	Excess								400	5	BLOCKED
Acetophenone	HO01002		1						1000	6	15/30 mg Tenax (SKC 226-35-03)
NIOSH S347	HO01003		1						1000	7	Acid Silica Gel (SKC 226-10-06)
OSHA ID200	HO01005		1						1000	8	100/50 mg Carbon Bead (SKC 226-80)
	Excess								1000	9	BLOCKED
TO11	HO01006		1						1000	10	DNPH Silica Gel (SKC 226-119)
	Excess								1000	_	BLOCKED
	Moisture		1						500	12	TOTAL
	Excess								5000	13	Excess

## RESEARCH FOUNDRY HO - SERIES SAMPLE PLAN

Method	Sample #	Data	Sample	Duplicate	Blank	Breakthrough	Spike	Spike Duplicate	Flow (ml/min)		Comments: System blank with core blower burners on. There was some sand in the hopper.				
4/26/2007															
THC, CH4, CO, CO <sub>2</sub> & NOx	HO011	X									TOTAL				
TO-17	HO01101		1						200	1	Carbopak charcoal				
	Excess								200	2	BLOCKED				
	Excess								200	3	BLOCKED				
	Excess								400	4	BLOCKED				
	Excess								400	5	BLOCKED				
Acetophenone	HO01102		1						1000	6	15/30 mg Tenax (SKC 226-35-03)				
NIOSH S347	HO01103		1						1000	7	Acid Silica Gel (SKC 226-10-06)				
OSHA ID200	HO01105		1						1000	8	100/50 mg Carbon Bead (SKC 226-80)				
	Excess								1000	9	BLOCKED				
TO11	HO01106		1						1000	10	DNPH Silica Gel (SKC 226-119)				
	Excess								1000	11	BLOCKED				
	Moisture		1						500	12	TOTAL				
	Excess								5000	13	Excess				

## **RESEARCH FOUNDRY HO - SERIES SAMPLE PLAN**

Method	Sample #	Data	Sample	Duplicate	Blank	Breakthrough	Spike	Spike Duplicate	Flow (ml/min)	Comments: System blank with core blower burners on and no sand in the hopper. No tube samples taken.
4/30/2007										
THC, CH4, CO, CO <sub>2</sub> & NOx	HO012	Х								TOTAL

Method	Sample #	Data	Sample	Duplicate	Blank	Breakthrough	Spike	Spike Duplicate	Flow (ml/min)	Train Channel	Comments:
4/25/2007											
THC, CH4, CO, CO <sub>2</sub> & NOx	HO001	Х									TOTAL
TO-17	HO00101		1						200	1	Carbopak charcoal
TO-17	HO00102				1				0		Carbopak charcoal
	Excess								200	2	BLOCKED
	Excess								200	3	BLOCKED
	Excess								400	4	BLOCKED
	Excess								400	5	BLOCKED
Acetophenone	HO00103		1						1000	6	15/30 mg Tenax (SKC 226-35-03)
Acetophenone	HO00104				1				0		15/30 mg Tenax (SKC 226-35-03)
NIOSH S347	HO00105		1						1000	7	Acid Silica Gel (SKC 226-10-06)
NIOSH S347	HO00106				1				0		Acid Silica Gel (SKC 226-10-06)
OSHA ID200			1						1000	8	100/50 mg Carbon Bead (SKC 226-80)
OSHA ID200	HO00110				1				0		100/50 mg Carbon Bead (SKC 226-80)
TO11			1						1000		DNPH Silica Gel (SKC 226-119)
TO11				1					1000	10	DNPH Silica Gel (SKC 226-119)
TO11					1				0		DNPH Silica Gel (SKC 226-119)
	Excess								1000		BLOCKED
	Moisture		1						500	12	TOTAL
	Excess								5000	13	Excess

#### **RESEARCH FOUNDRY HO - SERIES SAMPLE PLAN**

RESEARCH FOUNDRY HO - SERIES SAMPLE PLAN												
Method	Sample#	Data	Sample	Duplicate	Blank	Breakthrough	Spike	Spike Duplicate	Flow (ml/min)	Train Channel	Comments:	
4/25/2007												
THC, CH4, CO, CO <sub>2</sub> & NOx	HO002	Х									TOTAL	
TO-17	HO00201		1						200	1	Carbopak charcoal	
TO-17	HO00202			1					200	2	Carbopak charcoal	
	Excess								200	3	BLOCKED	
	Excess								400	4	BLOCKED	
	Excess								400	5	BLOCKED	
Acetophenone	HO00203		1						1000	6	15/30 mg Tenax (SKC 226-35-03)	
Acetophenone	HO00204			1					1000	7	15/30 mg Tenax (SKC 226-35-03)	
NIOSH S347	HO00205		1						1000	8	Acid Silica Gel (SKC 226-10-06)	
NIOSH S347	HO00206			1					1000	9	Acid Silica Gel (SKC 226-10-06)	
OSHA ID200	HO00208		1						1000	10	100/50 mg Carbon Bead (SKC 226-80)	
TO11	HO00209		1						1000	11	DNPH Silica Gel (SKC 226-119)	
	Moisture		1						500	12	TOTAL	
	Excess								5000	13	Excess	

#### RESEARCH FOUNDRY HO - SERIES SAMPLE PLAN

RESEARCH FOUNDRY HO - SERIES SAMPLE PLAN													
Method	Sample #	Data	Sample	Duplicate	Blank	Breakthrough	Spike	Spike Duplicate	Flow (ml/min)	Train Channel	Comments:		
4/26/2007													
THC, CH4, CO, CO <sub>2</sub> & NOx	HO003	Х									TOTAL		
TO-17	HO00301		1						200	1	Carbopak charcoal		
TO-17 MS	HO00302		1						200	2	Carbopak charcoal		
TO-17 MS	HO00303			1					200	3	Carbopak charcoal		
	Excess								400	4	BLOCKED		
	Excess								400	5	BLOCKED		
Acetophenone	HO00304		1						1000	6	15/30 mg Tenax (SKC 226-35-03)		
NIOSH S347	HO00305		1						1000	7	Acid Silica Gel (SKC 226-10-06)		
OSHA ID200	HO00308		1						1000	8	100/50 mg Carbon Bead (SKC 226-80)		
OSHA ID200	HO00309			1					1000	9	100/50 mg Carbon Bead (SKC 226-80)		
TO11	HO00310		1						1000	10	DNPH Silica Gel (SKC 226-119)		
	Excess								1000	11	BLOCKED		
	Moisture		1						500	12	TOTAL		
	Excess								5000	13	Excess		

<b>RESEARCH FOUNDRY HO -</b>	SERIES SAMPLE PLAN
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Method 4/26/2007	Sample #	Data	Sample	Duplicate	Blank	Breakthrough	Spike	Spike Duplicate	Flow (ml/min)	Train Channel	Comments:
4/20/2007											
THC, CH4, CO, CO <sub>2</sub> & NOx	HO004	Х									TOTAL
TO-17	HO00401		1						200	1	Carbopak charcoal
TO-17	HO00402					1			200	1	Carbopak charcoal
	Excess								200	2	BLOCKED
	Excess								200	3	BLOCKED
	Excess								400	4	BLOCKED
	Excess								400	5	BLOCKED
Acetophenone	HO00403		1						1000	6	15/30 mg Tenax (SKC 226-35-03)
NIOSH S347	HO00404		1						1000	7	Acid Silica Gel (SKC 226-10-06)
	Excess								1000	8	BLOCKED
OSHA ID200	HO00405		1						1000	9	100/50 mg Carbon Bead (SKC 226-80)
	Excess								1000	10	BLOCKED
TO11	HO00406		1						1000	11	DNPH Silica Gel (SKC 226-119)
	Moisture		1						500	12	BLOCKED
	Excess		1						5000	13	Excess

## **RESEARCH FOUNDRY HO - SERIES SAMPLE PLAN**

RESEARCH FOUNDRY NO - SERIES SAMPLE PLAN												
Method	Sample #	Data	Sample	Duplicate	Blank	Breakthrough	Spike	Spike Duplicate	Flow (ml/min)	Train Channel	Comments:	
4/26/2007												
THC, CH4, CO, CO <sub>2</sub> & NOx	HO005	Х									TOTAL	
TO-17	HO00501		1						200	1	Carbopak charcoal	
	Excess								200	2	BLOCKED	
	Excess								200	3	BLOCKED	
	Excess								400	4	BLOCKED	
	Excess								400	5	BLOCKED	
Acetophenone	HO00502		1						1000	6	15/30 mg Tenax (SKC 226-35-03)	
NIOSH S347	HO00503		1						1000	7	Acid Silica Gel (SKC 226-10-06)	
OSHA ID200	HO00504		1						1000	8	100/50 mg Carbon Bead (SKC 226-80)	
	Excess								1000	9	BLOCKED	
TO11	HO00505		1						1000	10	DNPH Silica Gel (SKC 226-119)	
	Excess								1000	11	BLOCKED	
	Moisture		1						500	12	TOTAL	
	Excess								5000	13	Excess	

## **RESEARCH FOUNDRY HO - SERIES SAMPLE PLAN**

RESEARCH FOUNDRY HO - SERIES SAMIFEE FEAR											
Method	Sample #	Data	Sample	Duplicate	Blank	Breakthrough	Spike	Spike Duplicate	Flow (ml/min)		Comments:
4/26/2007											
THC, CH4, CO, CO <sub>2</sub> & NOx	HO006	Х									TOTAL
TO-17	HO00601		1						200	1	Carbopak charcoal
	Excess								200	2	BLOCKED
	Excess								200	3	BLOCKED
	Excess								400	4	BLOCKED
	Excess								400	5	BLOCKED
Acetophenone	HO00602		1						1000	6	15/30 mg Tenax (SKC 226-35-03)
NIOSH S347	HO00603		1						1000	7	Acid Silica Gel (SKC 226-10-06)
OSHA ID200	HO00604		1						1000	8	100/50 mg Carbon Bead (SKC 226-80)
	Excess								1000	9	BLOCKED
TO11	HO00605								1000	10	DNPH Silica Gel (SKC 226-119)
	Excess								1000	11	BLOCKED
	Moisture		1						500	12	TOTAL
	Excess								5000	13	Excess

## <u>Series – 1413-122 HN</u>

# Core Making and Storage Fairmount Minerals Technisand XC30 Shell Sand Process Instructions

## **A.** The Experiment:

1. Measure total selected emissions from the making of step cores with Technisand XC30. This will also include a storage time.

#### **B.** Materials:

- 1. Technisand XC30, 3% resin (BOS).
- 2. Ashland Zipslip 109 W parting spray.
- **3.** Dow Corning #7 release agent.

## C. Machine setup.

- 1. Install and align the upgraded step core box onto the shell core machine.
- 2. Adjust gas and allow time for the east side of the shell box to reach 450F.
- **D.** Core Making test: Six (6) tests each having fifteen (15) approximately two (2) minute core cycles, with a thirty (30) minute storage cycle.

Note: It may take up to two (2) hours to temperature stabilize all the equipment.

Note: Use a smoke generator to verify that containment is achieved particularly around the shell box during machine use.

- 1. Prepare the core machine emission enclosure.
- 2. Cycle the core machine for 2-5 cycles or until background emissions are stable based on the THC and good core manufacture is achieved. Remove these cores from the enclosure.
- 3. When everybody is ready, start the emission sampling clock and open the sample train. Sample continuously for 15 to 20 core cycles, approximately thirty (30) minutes and while cores are in the hood, sample for another 30 minutes for a total of 60 minutes.
- **4.** Make the first core and place a thermocouple in it.
- **5.** Make cores continuously and place each core on the cart/table in the enclosure after removed from the machine. Any stoppage will impact the fugitive emission level.
- **6.** Record sample number, box temperature, and start time made of each core throughout the test.

Note: Occasionally, an incompletely cured core will occur. This does not necessarily constitute a failed sample emission cycle. Inferior cores should not be used for other purposes. Cores made for storage testing should be complete and no less than 0.1 pounds lighter than the median core weight for the group. Cores made for other testing may have the tip missing (0.05 pounds).

- 7. Do not stop making cores unless emission background level is allowed to re-stabilize before another test is begun.
- 8. Set up the sample train again and repeat the test for another fifteen core test. A total of six (6) tests are to be performed.
- **9.** Empty and clean the core machine and core sand mixer.

Tom Fennell Process Engineer

## <u>Series – 1413-115 HO</u>

## Core Making and Storage Fairmount Minerals Technisand 630BN Gold Shell Sand Process Instructions

## **A.** The Experiment:

1. Measure total selected emissions from the making of step cores with Technisand 630BN Gold. This will also include a storage time.

#### **B.** Materials:

- 1. Technisand 630BN Gold, 3% resin (BOS).
- **2.** Ashland Zipslip 109 W parting spray.
- **3.** Dow Corning #7 release agent.

## C. Machine setup.

- 1. Install and align the upgraded step core box onto the shell core machine.
- 2. Adjust gas and allow time for the east side of the shell box to reach 450F.
- **D.** Core Making test: Six (6) tests each having fifteen (15) approximately two (2) minute core cycles, with a thirty (30) minute storage cycle.

Note: It may take up to two (2) hours to temperature stabilize all the equipment.

Note: Use a smoke generator to verify that containment is achieved particularly around the shell box during machine use.

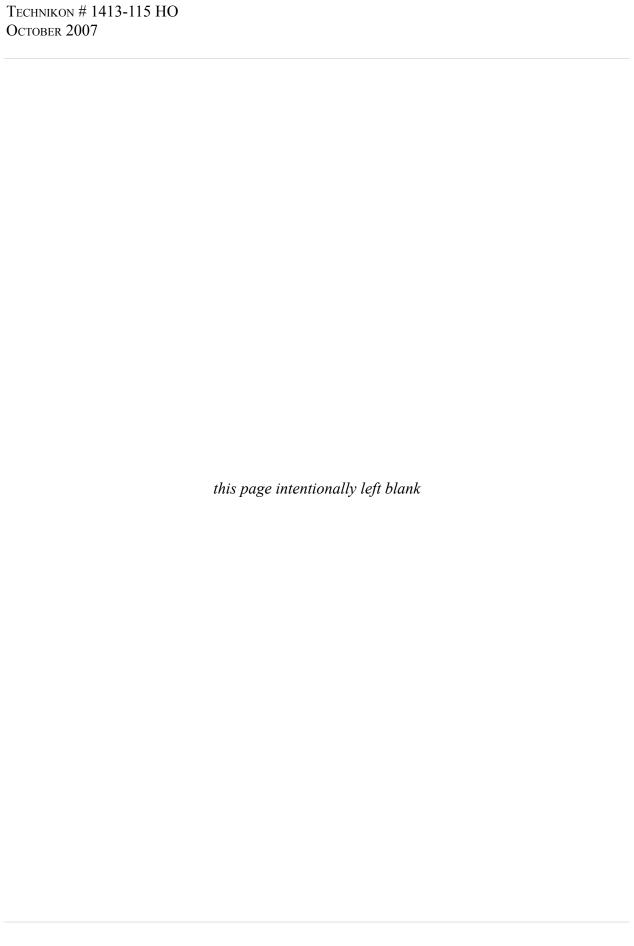
- 1. Prepare the core machine emission enclosure.
- 2. Cycle the core machine for 2-5 cycles or until background emissions are stable based on the THC and good core manufacture is achieved. Remove these cores from the enclosure.
- 3. When everybody is ready, start the emission sampling clock and open the sample train. Sample continuously for 15 to 20 core cycles, approximately thirty (30) minutes and while cores are in the hood, sample for another 30 minutes for a total of 60 minutes.
- **4.** Make the first core and place a thermocouple in it.
- **5.** Make cores continuously and place each core on the cart/table in the enclosure after removed from the machine. Any stoppage will impact the fugitive emission level.
- **6.** Record sample number, box temperature, and start time made of each core throughout the test.

Note: Occasionally, an incompletely cured core will occur. This does not necessarily constitute a failed sample emission cycle. Inferior cores should not be used for other purposes. Cores made for storage testing should be complete and no less than 0.1 pounds lighter than the median core weight for the group. Cores made for other testing may have the tip missing (0.05 pounds).

- 7. Do not stop making cores unless emission back ground level is allowed to re-stabilize before another test is begun.
- 8. Set up the sample train again and repeat the test for another fifteen core test. A total of six (6) tests are to be performed.
- **9.** Empty and clean the core machine and core sand mixer.

Tom Fennell Process Engineer

## APPENDIX B DETAILED EMISSION RESULTS AND QUANTITATION LIMITS



Emissions Core Make/Store and Storage, Test HN, lb/ton core

_			Emission	S Core IVI	ake/Store	and Sto	rage, res	L HIN, ID/L	/ton core		
ارا	POM	ΑF									Standard
₹	Ы	主		HN001	HN002	HN003	HN004	HN005	HN006	Average	Deviation
لِـــا		L	Test Dates	23-Apr-07	23-Apr-07	24-Apr-07	24-Apr-07	24-Apr-07	25-Apr-07		
Emi	ssio	n Inc	licators					· ·			
Ш			THC as Propane	5.17E-01	l	3.70E-01	4.59E-01	5.04E-01	8.22E-01	5.34E-01	1.71E-01
Ш			Non-Methane Hydrocarbons	ļ	ı	2.62E-01	2.64E-01	2.76E-01	5.28E-01	3.28E-01	1.30E-01
Ш			Sum of Target Analytes	5.61E-01	5.90E-01	5.60E-01	5.86E-01	6.10E-01	5.70E-01	5.79E-01	1.94E-02
Ш			Sum of Target HAPs	1.16E-01	1.32E-01	1.23E-01	1.29E-01	1.33E-01	1.14E-01	1.25E-01	8.05E-03
Ш			Sum of Target POMs	ND	ND	ND	ND	ND	ND	ND	NA
	cted		get HAPs and POMs								
TA		Н	Phenol	1.04E-01	1.19E-01	1.10E-01	1.16E-01	1.21E-01	1.03E-01	1.12E-01	7.44E-03
TA			Formaldehyde	1.06E-02	1.19E-02	1.18E-02	1.06E-02	1.14E-02	1.03E-02	1.11E-02	6.88E-04
TA		Н	Toluene	4.43E-04	≤PQL	≤PQL	1.35E-03	5.15E-04	≤PQL	4.46E-04	4.77E-04
TA		Н	Propionaldehyde (Propanal)	3.78E-04	4.41E-04	4.23E-04	3.62E-04	4.29E-04	5.45E-04	4.30E-04	6.44E-05
TA		Н	Acetaldehyde	3.84E-04	3.51E-04	3.20E-04	3.97E-04	4.29E-04	2.94E-04	3.62E-04	5.03E-05
TA		Н	Acrolein	I	1.23E-04	≤PQL	1.19E-04	1.19E-04	≤PQL	1.17E-04	5.22E-06
TA	Р	Н	Acenaphthalene	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA	Р	Н	Anthracene	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA	Р	Н	Dimethylnaphthalene, 1,2-	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA	Р	Н	Dimethylnaphthalene, 1,3-	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA	Р	Н	Dimethylnaphthalene, 1,5-	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA	Р	Н	Dimethylnaphthalene, 1,6-	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA	Р	Н	Dimethylnaphthalene, 1,8-	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA	Р	Н	Dimethylnaphthalene, 2,3-	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA	Р	Н	Dimethylnaphthalene, 2,6-	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA	Р	Н	Dimethylnaphthalene, 2,7-	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA	Р	Н	Methylnaphthalene, 1-	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA	Р	Н	Methylnaphthalene, 2-	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA	Р	Н	Naphthalene	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA	Р	Н	Trimethylnaphthalene, 2,3,5-	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA		Н	Benzene	≤PQL		≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA		Н	Biphenyl	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA		Н	Cresol, mp-	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA		Н	Cresol, o-	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA		Н	Ethylbenzene	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA		Н	Hexane	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA		Н	Styrene	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA		Н	Xylene, mp-	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA		Н	Xylene, o-	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA			Acetophenone	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
Add	ition	al S	elected Target Analytes								
TA			Ammonia	4.43E-01	4.56E-01	4.36E-01	4.56E-01	4.75E-01	4.54E-01	4.53E-01	1.34E-02
TA			Crotonaldehyde	5.15E-04	4.71E-04	4.38E-04	4.16E-04	4.77E-04	5.72E-04	4.81E-04	5.61E-05
TA			2-Butanone (MEK)	3.82E-04	2.87E-04	3.32E-04	3.23E-04	3.68E-04	3.15E-04	3.35E-04	3.50E-05
TA			Benzaldehyde	≤PQL	2.12E-04	2.89E-04	2.24E-04	2.62E-04	3.84E-04	2.47E-04	9.04E-05
TA			Butyraldehyde/Methacrolein	3.13E-04	2.25E-04	2.36E-04	2.04E-04	2.33E-04	2.62E-04	2.46E-04	3.81E-05
TA			Pentanal (Valeraldehyde)	I	≤PQL	≤PQL	≤PQL	1.14E-04	1.27E-04	1.15E-04	6.71E-06
TA			Cyclohexane	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
					•	•		•	Continue	d on nex	(Ansart

(continued on next page)

I=invalidated

NA=not applicable ≤PQL=less than or equal to pratical quantation limit Emissions Core Make/Store and Storage, Test HN, lb/ton core (continued)

February   February				Emission	3 OUIC IVI	akerotore	and old	rage, res	11111, 115/1		Continue	
Test Dates	4	등	ΑP		LINIOOA	LINIOOO	LINIOOO	1111004	LINIOOE	LINIOOC	A	Standard
TA	È	ے	Ì	Total Dates							Average	Deviation
TA	- A	_									- -	
TA		_										
TA												
TA												
TA				71 /								
TA												
TA												
TA         Hexaldehyde         ≤PQL         SPQL         ≤PQL         SPQL												
TA         Indan         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         SPQL         ≤PQL         SPQL         SPQL <th< th=""><th></th><th></th><th></th><th></th><th></th><th></th><th></th><th></th><th></th><th></th><th></th><th></th></th<>												
TA         Nonane         SPQL         SPQL <t< th=""><th></th><th></th><th></th><th></th><th></th><th></th><th></th><th></th><th></th><th></th><th></th><th></th></t<>												
TA         □ o,m.pTolualdehyde         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         SPQL         SPQL<				Indan	≤PQL		≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA         Octane         ≤PQL         SPQL         ≤PQL         SPQL         SPQL <t< th=""><th>TA</th><th></th><th></th><th></th><th>≤PQL</th><th>≤PQL</th><th>≤PQL</th><th>≤PQL</th><th>≤PQL</th><th>≤PQL</th><th>≤PQL</th><th>NA</th></t<>	TA				≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA         Propylbenzene, n-         ≤PQL         ≤PQL <pql< th="">         ≤PQL         <pql< th=""> <pql< th=""></pql<></pql<></pql<>	TA			o,m,p-Tolualdehyde	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA         Tetradecane         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         SPQL							≤PQL	≤PQL		≤PQL	≤PQL	NA
TA         Trimethylbenzene, 1,2,3-         ≤PQL         ≤PQL <pql< th=""> <pql< th="">         SPQL         SP</pql<></pql<>				Propylbenzene, n-	≤PQL		≤PQL	≤PQL	≤PQL		≤PQL	NA
TA         Trimethylbenzene, 1,2,4-         ≤PQL         ≤P	TA			Tetradecane				≤PQL				NA
TA         Inimethylbenzene, 1,3,5-         ≤PQL         SPQL         ≤PQL         ≤PQL         SPQL         SP	TA			Trimethylbenzene, 1,2,3-	≤PQL		≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA         Undecane         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         NA           TA         alpha-Methylstyrene         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         SPQL         SPQL         SPQL         NA           TA         Butylbenzene, n-         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         SPQL         SPQL         SPQL         NA           TA         Butylbenzene, sec-         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         SPQL         SPQL         SPQL         NA           TA         Butylbenzene, tert-         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         SPQL         SPQL         SPQL         NA           TA         Diethylbenzene, tert-         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         SPQL         SPQL         SPQL         NA           TA         Diethylbenzene, t.2-         ≤PQL         ≤PQL         ≤PQL         ≤PQL         SPQL         SPQL         SPQL         NA           TA         Diisethylbenzene, t.3-         ≤PQL         ≤PQL	TA			Trimethylbenzene, 1,2,4-	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA         alpha-Methylstyrene         ≤PQL         ≤PQL <th>TA</th> <th></th> <th></th> <th>Trimethylbenzene, 1,3,5-</th> <th>≤PQL</th> <th>≤PQL</th> <th>≤PQL</th> <th>≤PQL</th> <th>≤PQL</th> <th>≤PQL</th> <th>≤PQL</th> <th>NA</th>	TA			Trimethylbenzene, 1,3,5-	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA         Butylbenzene, n-         ≤PQL	TA			Undecane	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA         Butylbenzene, sec-         ≤PQL         ≤PQL <pql< th="">         ≤PQL         <pql< th=""> <pql< th="" th<=""><th>TA</th><th></th><th></th><th>alpha-Methylstyrene</th><th>≤PQL</th><th>≤PQL</th><th>≤PQL</th><th>≤PQL</th><th>≤PQL</th><th>≤PQL</th><th>≤PQL</th><th>NA</th></pql<></pql<></pql<></pql<></pql<></pql<></pql<></pql<></pql<></pql<></pql<></pql<></pql<></pql<></pql<></pql<></pql<></pql<></pql<></pql<></pql<></pql<></pql<></pql<></pql<></pql<></pql<></pql<></pql<></pql<>	TA			alpha-Methylstyrene	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA         Butylbenzene, tert-         ≤PQL         ≤PQL <th>TA</th> <th></th> <th></th> <th>Butylbenzene, n-</th> <th>≤PQL</th> <th>≤PQL</th> <th>≤PQL</th> <th>≤PQL</th> <th>≤PQL</th> <th>≤PQL</th> <th>≤PQL</th> <th>NA</th>	TA			Butylbenzene, n-	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA         Cymene, p-         ≤PQL         ≤PQL <pql< th=""> <pql< th="">         ≤PQL         <pql< th=""></pql<></pql<></pql<>	TA			Butylbenzene, sec-	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA         Cymene, p-         ≤PQL         ≤PQL <pql< th=""> <pql< th="">         ≤PQL         <pql< th=""></pql<></pql<></pql<>	TA			Butylbenzene, tert-	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA         Diethylbenzene, 1,4-         ≤PQL         ≤PQL </th <th>TA</th> <th></th> <th></th> <th></th> <th>≤PQL</th> <th>≤PQL</th> <th>≤PQL</th> <th>≤PQL</th> <th>≤PQL</th> <th>≤PQL</th> <th>≤PQL</th> <th>NA</th>	TA				≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA         Diisopropylbenzene, 1,3-         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         NA           TA         Dimethylphenol, 2,3-         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         NA           TA         Dimethylphenol, 2,5-         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         SPQL         NA           TA         Dimethylphenol, 3,4-         ≤PQL         ≤PQL         ≤PQL         ≤PQL         SPQL         SPQL         SPQL         SPQL         SPQL         NA           TA         Dimethylphenol, 3,5-         ≤PQL         ≤PQL         ≤PQL         ≤PQL         SPQL         SPQL         SPQL         SPQL         NA           TA         Dimethylphenol, 3,5-         ≤PQL         ≤PQL         ≤PQL         ≤PQL         SPQL         SPQL         SPQL         NA           TA         Indene         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         SPQL </th <th>TA</th> <th></th> <th></th> <th>Diethylbenzene, 1,2-</th> <th>≤PQL</th> <th>≤PQL</th> <th>≤PQL</th> <th>≤PQL</th> <th>≤PQL</th> <th>≤PQL</th> <th>≤PQL</th> <th>NA</th>	TA			Diethylbenzene, 1,2-	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA         Dimethylphenol, 2,3-         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         NA           TA         Dimethylphenol, 2,5-         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         SPQL         SPQL         SPQL         NA           TA         Dimethylphenol, 3,4-         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         SPQL         SPQL         SPQL         NA           TA         Dimethylphenol, 3,5-         ≤PQL         ≤PQL         ≤PQL         ≤PQL         SPQL         SPQL         SPQL         NA           TA         Dimethylphenol, 3,5-         ≤PQL         ≤PQL         ≤PQL         ≤PQL         SPQL         SPQL         SPQL         NA           TA         Dimethylphenol, 3,5-         ≤PQL         ≤PQL         ≤PQL         ≤PQL         SPQL         SPQL         SPQL         NA           TA         Indene         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         SPQL         SPQL         NA           TA         Indecane         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL	TA			Diethylbenzene, 1,4-	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA         Dimethylphenol, 2,5-         ≤PQL         NA           TA         Dimethylphenol, 3,4-         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         NA           TA         Dimethylphenol, 3,5-         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         SPQL         SPQL         SPQL         NA           TA         Ethyltoluene, 4-         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         SPQL         SPQL         SPQL         NA           TA         Indene         ≤PQL         ≤PQL         ≤PQL         ≤PQL         SPQL         SPQL         NA           TA         Isobutylbenzene         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         NA           TA         Isopropylbenzene         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         NA           TA         Tridecane         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQ	TA			Diisopropylbenzene, 1,3-	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA         Dimethylphenol, 3,4-         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         NA           TA         Dimethylphenol, 3,5-         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         SPQL         <	TA			Dimethylphenol, 2,3-	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA         Dimethylphenol, 3,5-         ≤PQL         NA           TA         Indene         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ×PQL	TA			Dimethylphenol, 2,5-	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA         Ethyltoluene, 4-         ≤PQL         NA           TA         Indene         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         NA           TA         Isopropylbenzene         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         SPQL         SPQL         SPQL         SPQL         SPQL         SPQL         SPQL         NA           TA         Trimethylphenol, 2,3,5-         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         SPQL         SPQL         NA           TA         Trimethylphenol, 2,4,6-         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         SPQL         SPQL         NA           0.00E+00         Methane         I         I         1.08E-01         1.95E-01         2.28E-01         2.94E-01         2.06E-01         7.74E-02           Carbon Monoxide         1.16E-01         1.26E-01         1.59E-01         1.04E-01         1.04E-01         1.44E-01         1.26E-01         2.15E-02	TA			Dimethylphenol, 3,4-	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA         Indene         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         NA           TA         Isobutylbenzene         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         NA           TA         Isopropylbenzene         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         SPQL         SPQL         SPQL         SPQL         NA           TA         Trimethylphenol, 2,3,5-         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         SPQL         SPQL         SPQL         NA           TA         Trimethylphenol, 2,4,6-         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         SPQL         NA           0.00E+00         Methane         I         I         1.08E-01         1.95E-01         2.28E-01         2.94E-01         2.06E-01         7.74E-02           Carbon Monoxide         1.16E-01         1.26E-01         1.59E-01         1.04E-01         1.04E-01         1.44E-01         1.26E-01         2.15E-02           Nitrogen Oxides         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL <th>TA</th> <th></th> <th></th> <th>Dimethylphenol, 3,5-</th> <th>≤PQL</th> <th>≤PQL</th> <th>≤PQL</th> <th>≤PQL</th> <th>≤PQL</th> <th>≤PQL</th> <th>≤PQL</th> <th>NA</th>	TA			Dimethylphenol, 3,5-	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA         Indene         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         NA           TA         Isobutylbenzene         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         NA           TA         Isopropylbenzene         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         SPQL         SPQL         SPQL         SPQL         NA           TA         Trimethylphenol, 2,3,5-         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         SPQL         SPQL         SPQL         NA           TA         Trimethylphenol, 2,4,6-         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         SPQL         NA           0.00E+00         Methane         I         I         1.08E-01         1.95E-01         2.28E-01         2.94E-01         2.06E-01         7.74E-02           Carbon Monoxide         1.16E-01         1.26E-01         1.59E-01         1.04E-01         1.04E-01         1.44E-01         1.26E-01         2.15E-02           Nitrogen Oxides         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL <th>TA</th> <th></th> <th></th> <th>Ethyltoluene, 4-</th> <th>≤PQL</th> <th>≤PQL</th> <th>≤PQL</th> <th>≤PQL</th> <th>≤PQL</th> <th>≤PQL</th> <th>≤PQL</th> <th>NA</th>	TA			Ethyltoluene, 4-	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA         Isobutylbenzene         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         NA           TA         Isopropylbenzene         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         SPQL         SPQL         SPQL         NA           TA         Trimethylphenol, 2,3,5-         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         SPQL         SPQL         SPQL         NA           TA         Trimethylphenol, 2,4,6-         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         SPQL         NA           0.00E+00         Methane         I         I         1.08E-01         1.95E-01         2.28E-01         2.94E-01         2.06E-01         7.74E-02           Carbon Monoxide         1.16E-01         1.26E-01         1.59E-01         1.04E-01         1.08E-01         1.44E-01         1.26E-01         2.15E-02           Nitrogen Oxides         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL <pql< th=""> <p< th=""><th></th><th></th><th></th><th></th><th></th><th></th><th></th><th></th><th></th><th></th><th></th><th>NA</th></p<></pql<></pql<></pql<></pql<></pql<></pql<></pql<></pql<>												NA
TA         Isopropylbenzene         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         NA           TA         Tridecane         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         NA           TA         Trimethylphenol, 2,4,6-         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         NA           TA         Methane         I         I         1.08E-01         1.95E-01         2.28E-01         2.94E-01         2.06E-01         7.74E-02           Carbon Monoxide         1.16E-01         1.26E-01         1.59E-01         1.04E-01         1.04E-01         1.44E-01         1.26E-01         2.15E-02           Nitrogen Oxides         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         NA	TA			Isobutylbenzene	≤PQL	≤PQL	≤PQL		≤PQL	≤PQL	≤PQL	NA
TA         Trimethylphenol, 2,3,5-         ≤PQL         NA           Timethylphenol, 2,4,6-         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         NA           0.00E+00           Methane         I         I         1.08E-01         1.95E-01         2.28E-01         2.94E-01         2.06E-01         7.74E-02           Carbon Monoxide         1.16E-01         1.26E-01         1.59E-01         1.04E-01         1.08E-01         1.44E-01         1.26E-01         2.15E-02           Nitrogen Oxides         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         >PQL         >PQL         NA	TA			Isopropylbenzene	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA         Trimethylphenol, 2,4,6-         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         NA           0.00E+00           Methane         I         I         1.08E-01         1.95E-01         2.28E-01         2.94E-01         2.06E-01         7.74E-02           Carbon Monoxide         1.16E-01         1.26E-01         1.59E-01         1.04E-01         1.04E-01         1.44E-01         1.26E-01         2.15E-02           Nitrogen Oxides         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         NA	TA			Tridecane	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
0.00E+00           Methane         I         I         1.08E-01         1.95E-01         2.28E-01         2.94E-01         2.06E-01         7.74E-02           Carbon Monoxide         1.16E-01         1.26E-01         1.59E-01         1.08E-01         1.08E-01         1.44E-01         1.26E-01         2.15E-02           Nitrogen Oxides         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         NA	TA			Trimethylphenol, 2,3,5-	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
0.00E+00           Methane         I         I         1.08E-01         1.95E-01         2.28E-01         2.94E-01         2.06E-01         7.74E-02           Carbon Monoxide         1.16E-01         1.26E-01         1.59E-01         1.04E-01         1.08E-01         1.44E-01         1.26E-01         2.15E-02           Nitrogen Oxides         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         NA	TA					≤PQL		≤PQL		≤PQL		NA
Methane         I         I         1.08E-01         1.95E-01         2.28E-01         2.94E-01         2.06E-01         7.74E-02           Carbon Monoxide         1.16E-01         1.26E-01         1.59E-01         1.04E-01         1.08E-01         1.44E-01         1.26E-01         2.15E-02           Nitrogen Oxides         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         NA	0.00	E+00	)	71								
Carbon Monoxide         1.16E-01         1.26E-01         1.59E-01         1.04E-01         1.08E-01         1.44E-01         1.26E-01         2.15E-02           Nitrogen Oxides         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         ≤PQL         NA				Methane	I		1.08E-01	1.95E-01	2.28E-01	2.94E-01	2.06E-01	7.74E-02
Nitrogen Oxides ≤PQL ≤PQL ≤PQL ≤PQL ≤PQL ≤PQL ≤PQL NA					1.16E-01	1.26E-01						
				Sulfur Dioxide	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
0.00E+00												

I=invalidated NA=not applicable

≤PQL=less than or equal to pratical quantation limit

Detailed Emissions Core Make/Store and Storage, Test HN, lb/lb binder

			Detailed Emiss	SIONS CO	e Wake/S	tore and	Storage,	Test min,	וווט טווט	Jei	
≱	POM	₽		1111004	LINIOOO	LINIOOS	LINIOOA	LINIOGE	LINIOOC	A	Standard
브	Ā	ヹ	T (D)	HN001	HN002	HN003	HN004	HN005	HN006	Average	Deviation
<u> </u>	<u> </u>	Ļ	Test Dates	23-Apr-07	23-Apr-07	24-Apr-07	24-Apr-07	24-Apr-07	25-Apr-07		
Emi	SSIOI	n Inc	licators	0.005.00		0045.00	7.005.00	0.005.00	4.445.00	0.40=.00	0.005.00
$\vdash$			THC as Propane	8.88E-03	<u> </u>	6.34E-03	7.86E-03	8.66E-03	1.41E-02	9.16E-03	2.92E-03
$\vdash$			Non-Methane Hydrocarbons	0.005.00	1 0 1 5 0 0	4.49E-03	4.53E-03	4.74E-03	9.04E-03	5.63E-03	2.23E-03
$\vdash$			Sum of Target Analytes	9.62E-03	1.01E-02	9.72E-03	1.03E-02	1.07E-02	9.80E-03	1.00E-02	4.22E-04
Ш			Sum of Target HAPs	1.99E-03	2.31E-03	2.13E-03	2.26E-03	2.35E-03	1.97E-03	2.17E-03	1.63E-04
		<u> </u>	Sum of Target POMs	ND	ND	ND	ND	ND	ND	ND	NA
-	cted	_	get HAPs and POMs	4 705 00	0.005.00	404500	0.005.00	0.405.00	4 775 00	4.055.00	4.545.04
TA		Н	Phenol	1.79E-03	2.08E-03	1.91E-03	2.03E-03	2.12E-03	1.77E-03	1.95E-03	1.51E-04
TA		Н	Formaldehyde	1.82E-04	2.09E-04	2.05E-04	1.87E-04	2.00E-04	1.77E-04	1.93E-04	1.31E-05
TA		Н	Toluene	7.60E-06	≤PQL	≤PQL	2.37E-05	9.07E-06	≤PQL	7.79E-06	8.40E-06
TA		Н	Propionaldehyde (Propanal)	6.48E-06	7.74E-06	7.35E-06	6.37E-06	7.55E-06	9.38E-06	7.48E-06	1.09E-06
TA		Н	Acetaldehyde	6.59E-06	6.15E-06	5.55E-06	6.97E-06	7.55E-06	5.06E-06	6.31E-06	9.18E-07
TA	_	Н	Acrolein		2.16E-06	≤PQL	2.08E-06	2.09E-06	≤PQL	2.03E-06	1.15E-07
TA	Р	Н	Acenaphthalene	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA	Р	Н	Anthracene	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA	Р	Н	Dimethylnaphthalene, 1,2-	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA	Р	Н	Dimethylnaphthalene, 1,3-	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA	Р	Н	Dimethylnaphthalene, 1,5-	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA	Р	Н	Dimethylnaphthalene, 1,6-	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA	Р	Н	Dimethylnaphthalene, 1,8-	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA	Р	Н	Dimethylnaphthalene, 2,3-	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA	Р	Н	Dimethylnaphthalene, 2,6-	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA	Р	Н	Dimethylnaphthalene, 2,7-	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA	Р	Н	Methylnaphthalene, 1-	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA	Р	Н	Methylnaphthalene, 2-	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA	Р	Н	Naphthalene	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA	Р	Н	Trimethylnaphthalene, 2,3,5-	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA		Н	Benzene	≤PQL	ı	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA		Н	Biphenyl	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA		Н	Cresol, mp-	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA		Н	Cresol, o-	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA		Н	Ethylbenzene	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA		Н	Hexane	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA		Н	Styrene	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA		Н	Xylene, mp-	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA		Н	Xylene, o-	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA			Acetophenone	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
	ition	al S	elected Target Analytes								
TA			Ammonia	7.60E-03	7.75E-03	7.56E-03	8.02E-03	8.36E-03	7.80E-03	7.85E-03	2.99E-04
TA			Crotonaldehyde	8.83E-06	8.25E-06	7.60E-06	7.30E-06	8.40E-06	9.84E-06	8.37E-06	9.10E-07
TA			2-Butanone (MEK)	6.55E-06	5.03E-06	5.77E-06	5.68E-06	6.49E-06	5.41E-06	5.82E-06	5.97E-07
TA			Benzaldehyde	≤PQL	3.72E-06	5.02E-06	3.94E-06	4.62E-06	6.60E-06	4.30E-06	1.56E-06
TA			Butyraldehyde/Methacrolein	5.38E-06	3.95E-06	4.10E-06	3.58E-06	4.11E-06	4.51E-06	4.27E-06	6.19E-07
TA			Pentanal (Valeraldehyde)	1	≤PQL	≤PQL	≤PQL	2.00E-06	2.18E-06	1.98E-06	1.18E-07
TA			Cyclohexane	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA

(continued on next page)

I=invalidated
NA=not applicable
≤PQL=less than or equal to pratical quantation limit

Detailed Emissions Core Make/Store and Storage, Test HN, lb/lb binder (continued)

			Detailed Emis	310113 001	e make/c	tore and	otorage,	16311111,	ווע טווע	Jei (conti	
l≱	POM	HAP		HN001	HN002	HN003	HN004	HN005	HN006	Average	Standard Deviation
	_	_	Test Dates	23-Apr-07	23-Apr-07	24-Apr-07	24-Apr-07	24-Apr-07	25-Apr-07	_	_
TA			Decane	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA			Diethylbenzene, 1,3-	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA			Dimethylphenol, 2,4-	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA			Dimethylphenol, 2,6-	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA			Dodecane	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA			Ethyltoluene, 2-	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA			Ethyltoluene, 3-	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA			Heptane	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA			Hexaldehyde	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA			Indan	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA			Nonane	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA			o,m,p-Tolualdehyde	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA			Octane	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA			Propylbenzene, n-	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA			Tetradecane	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA			Trimethylbenzene, 1,2,3-	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA			Trimethylbenzene, 1,2,4-	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA			Trimethylbenzene, 1,3,5-	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA			Undecane	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA			alpha-Methylstyrene	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA			Butylbenzene, n-	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA			Butylbenzene, sec-	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA			Butylbenzene, tert-	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA			Cymene, p-	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA			Diethylbenzene, 1,2-	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA			Diethylbenzene, 1,4-	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA			Diisopropylbenzene, 1,3-	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA			Dimethylphenol, 2,3-	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA	_		Dimethylphenol, 2,5-	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA	_		Dimethylphenol, 3,4-	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA	_		Dimethylphenol, 3,5-	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA			Ethyltoluene, 4-	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA			Indene	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA			Isobutylbenzene	≤PQL	≤PQL ≤PQL	≤PQL	≤PQL ≤PQL	≤PQL ≤PQL	≤PQL	≤PQL	NA NA
TA			Isopropylbenzene	≤PQL	≤PQL ≤PQL	≤PQL			≤PQL	≤PQL	NA NA
TA			Tridecane	≤PQL		≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA NA
TA TA			Trimethylphenol, 2,3,5-	≤PQL ≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA NA
	040.0	C-:-	Trimethylphenol, 2,4,6-		≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	INA
Sele	ctea	Crit	eria Pollutants and Greenho Carbon Dioxide	2.71E-01	2.63E-01	2.83E-01	2.82E-01	2.75E-01	2.65E-01	2.73E-01	8.29E-03
			Methane	Z./ IE-UI	2.03E-01	2.83E-01 1.85E-03	3.34E-03	3.92E-03	5.04E-03	3.54E-03	1.33E-03
	$\dashv$		Carbon Monoxide	1.99E-03	2.16E-03	2.72E-03	1.79E-03	1.85E-03	2.47E-03	2.16E-03	3.68E-04
			Nitrogen Oxides	1.99E-03 ≤PQL	2.10E-03 ≤PQL	2.72E-03 ≤PQL	1.79E-03 ≤PQL	1.05E-03 ≤PQL	2.47E-03 ≤PQL	2.10E-03 ≤PQL	3.00E-04 NA
			Sulfur Dioxide	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA NA
			Journal Dioxide	≥rul	≥rul	≥rul	≥rul	≥rų∟	≥rul	≥rų∟	INA

I=invalidated NA=not applicable

≤PQL=less than or equal to pratical quantation limit

Practical Reporting Limits, Core Make/Store and Storage, Test

Practical Reporting Limits, Core Make/Store and Storage, Test HN, Ib/ton core

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Analyte         Ibitb Binder         Analyte         Ibitb Binder         Analyte         Ibitb Binder           2-Butanone (MEK)         1.91E-06         Acena         2.11E-06         Acena           Accentaphthalene         2.14E-06         Acenaphthalene         2.11E-06         Accentaphthalene           Accentaphthalene         2.14E-06         Accentaphthalene         2.11E-06         Accentaphthalene           Accentaphthalene         2.14E-06         Ethyllulenene, 3-         2.11E-06         Accentaphthalene           Anthracene         2.11E-06         Ethyllulenene, 3-         2.11E-06         Action           Anthracene         2.11E-06         Bernzade         2.11E-06         Bernzade           Buylbenzane, propertion         2.11E-06         Bernzade         2.11E-06         Bernzade           Buylbenzane, propertion         2.11E-06         Buylbenzane         2.11E-06         Buylbenzane           Carbon Monoxide </th <th></th> <th>MN, ID,</th> <th>HN, ID/ID BINGEr</th> <th></th> <th></th>		MN, ID,	HN, ID/ID BINGEr		
1.91E-06   Dodecane   2.11E-06   Ethylbenzene   2.11E-06   Ethylbenzene   2.11E-06   Ethylbenzene   2.11E-06   Ethylbenzene   2.11E-06   Ethylbenzene   2.11E-06   Ethylbenzene   2.11E-06   Ethylptoluene, 3	Analyte	lb/lb Binder	Analyte	lb/lb Binder	
191E-06   Ethyllotuene, 2-   2.11E-06   Ethyllotuene, 3-   2.11E-06   Ethyllotuene, 3-   2.11E-06   Ethyllotuene, 4-   2.11E-06   Hexane   2.11E-06   Hexane   2.11E-06   Hexane   2.11E-06   Hodan   2.	2-Butanone (MEK)	1.91E-06	Dodecane	2.11E-06	2-Bui
1.91E-06   Ethyltoluene, 2-	Acenaphthalene	2.11E-06	Ethylbenzene	2.11E-06	Acen
2.45E-04   Ethyltoluene, 3- 2.11E-06   Ethyltoluene, 4- 2.11E-06   Formaldehyde	Acetaldehyde	1.91E-06	Ethyltoluene, 2-	2.11E-06	Aceta
1.91E-06   Ethyltoluene, 4-   2.11E-06   Ethyltoluene, 4-   2.11E-06   Heptane   2.11E-06   Heptane   2.11E-06   Hexalchyde   2.11E-06   Hexalchyde   2.11E-06   Hexalchyde   2.11E-06   Hexalchyde   2.11E-06   Indan   Indan   2.11E-06   Indan   I	Acetophenone	2.45E-04	Ethyltoluene, 3-	2.11E-06	Acetc
2.11E-06         Formaldehyde         1.91E-06           1.11E-06         Hexandehyde         2.11E-06           1.91E-06         Hexandehyde         2.11E-06           2.11E-06         Indan         2.11E-06           2.11E-06         Indane         2.11E-06           2.11E-06         Indane         2.11E-06           Indene         2.11E-06	Acrolein	1.91E-06	Ethyltoluene, 4-	2.11E-06	Acrol
ee         7.41E-05         Heptane         2.11E-06           Hexaldehyde         1.91E-06           Hexaldehyde         1.91E-06           ene, ne, sec.         2.11E-06           ene, pe, rese, sec.         2.11E-06           ene, sec.         2.11E-06           ene, sec.         2.11E-06           ene, sec.         2.11E-06           lander         2.11E-06           lander         2.11E-06           lander         2.11E-06           lander         2.11E-06           lander         3.18E-06           lander         3.11E-06           lande	alpha-Methylstyrene	2.11E-06	Formaldehyde	1.91E-06	alpha
2.11E-06         Hexaldehyde         1.91E-06           1.91E-06         Hexane         2.11E-06           2.11E-06         Indene         2.11E-06           2.11E-06         Isobutylbenzene         2.11E-06           2.11E-06         Isobutylbenzene         2.11E-06           2.11E-06         Isobutylbenzene         2.11E-06           3.18E-06         Methylnaphthalene, 1-         2.11E-06           6.32E-04         Methylnaphthalene, 2-         2.11E-06           8.32E-04         Methylnaphthalene, 2-         2.11E-06           1.91E-06         Nonane         2.11E-06           1.91E-06         Octane         2.11E-06           2.11E-06         Pentanal (Valeraldehyde)         1.91E-06           2.11E-06         Porpylbenzene, n-         2.11E-06           Propylbenzene, n-         2.11E-06           Styrene         2.11E-06           Styrene         2.11E-06           Styrene         2.11E-06           Styrene         2.11E-06           Styrene         2.11E-06           Timethylbenzene, 1.2.3-         2.11E-06           Timethylbenzene, 1.2.4-         2.11E-06           Timethylbenzene, 1.2.3-         2.11E-06	Ammonia	7.41E-05	Heptane	2.11E-06	Amm
1.91E-06         Hexane         2.11E-06           2.11E-06         Indan         2.11E-06           2.11E-06         Isobulylbenzene         2.11E-06           2.11E-06         Methynaphthalene, 1-         2.11E-06           3.18E-06         Methynaphthalene, 2-         2.11E-06           6.32E-04         Methynaphthalene, 2-         2.11E-06           1.91E-06         Naphthalene, 2-         2.11E-06           2.11E-06         Nitrogen Oxides         2.11E-06           1.91E-06         O.m.p-Tolualdehyde         2.11E-06           2.11E-06         O.m.p-Tolualdehyde         2.11E-06           2.11E-06         O.m.p-Tolualdehyde         2.11E-06           2.11E-06         Phenol         2.11E-06           Phenol         2.11E-06         Phenol           2.11E-06         Propylbenzene, n-         2.11E-06           Phenol         Styrene         2.11E-06           Phenol         Propionaldehyde (Propanal)         1.91E-06           Phenol         Styrene         2.11E-06           Styrene         Timethylbenzene, n-         2.11E-06           Timethylbenzene, n-         2.11E-06           Timethylbenzene, n-         2.11E-06           Timethyl	Anthracene	2.11E-06	Hexaldehyde	1.91E-06	Anthi
2.11E-06         Inden         2.11E-06           2.11E-06         Indene         2.11E-06           2.11E-06         Isobutylbenzene         2.11E-06           2.11E-06         Methylnaphthalene, 1-         2.11E-06           3.18E-06         Methylnaphthalene, 2-         2.11E-06           6.32E-04         Methylnaphthalene, 2-         2.11E-06           6.32E-04         Methylnaphthalene, 2-         2.11E-06           1.91E-06         Nitrogen Oxides         2.11E-06           2.11E-06         Nonane         2.11E-06           2.11E-06         Octane         2.11E-06           2.11E-06         Pentanal (Valeraldehyde)         1.91E-06           2.11E-06         Pentanal (Valeraldehyde)         1.91E-06           2.11E-06         Popylbenzene, n-         2.11E-06           2.11E-06         Propylbenzene, n-         2.11E-06           2.11E-06         Propylbenzene, n-         2.11E-06           2.11E-06         Title-ofe           2.11E-06         Title-ofe           2.11E-06         Title-ofe           2.11E-06         Title-ofe           2.11E-06         Title-ofe           2.11E-06         Title-ofe           2.11E-06	Benzaldehyde	1.91E-06	Hexane	2.11E-06	Benz
2.11E-06         Indene         2.11E-06           2.11E-06         Isobutybenzene         2.11E-06           2.11E-06         Sporopybenzene         2.11E-06           3.18E-06         Methylnaphthalene, 1-         2.11E-06           6.32E-04         Methylnaphthalene, 2-         2.11E-06           6.32E-04         Methylnaphthalene, 2-         2.11E-06           A.02E-04         Nonane         2.11E-06           Nonane         2.11E-06         3.18E-06           Doctane         2.11E-06         3.11E-06           2.11E-06         Dentanal (Valeraldehyde)         2.11E-06           2.11E-06         Propionaldehyde (Propanal)         1.91E-06           2.11E-06         Propionaldehyde (Propanal)         2.11E-06           2.11E-06         Propylbenzene, n-         2.11E-06           2.11E-06         Sulfur Dioxide         2.33E-04           2.11E-06         Sulfur Dioxide         2.31E-06           2.11E-06         Timethylbenzene, 1.2.4         2.11E-06           2.11E-06         Timethylbenzene, 1.2.4         2.11E-06           2.11E-06         Trimethylbenzene, 1.2.4         2.11E-06           2.11E-06         Trimethylpenol, 2.3.5-         2.11E-06           2.11E-	Benzene	2.11E-06	Indan	2.11E-06	Benz
2.11E-06         Isobutylbenzene         2.11E-06           2.11E-06         Isopropylbenzene         2.11E-06           2.11E-06         Methylnaphthalene, 1-         2.11E-06           6.32E-04         Methylnaphthalene, 2-         2.11E-06           4.02E-04         Naphthalene         2.11E-06           2.11E-06         Norane         2.11E-06           2.11E-06         Octane         2.11E-06           2.11E-06         Phanol         2.11E-06           2.11E-06         Portanal (Valeraldehyde)         1.91E-06           2.11E-06         Propylbenzene, n-         2.11E-06           2.11E-06         Propylbenzene, n-         2.11E-06           2.11E-06         Styrene         2.11E-06           2.11E-06         Suffur Dioxide         2.53E-05           2.11E-06         Suffur Dioxide         2.11E-06           2.11E-06         Tetadecane         2.11E-06           2.11E-06         Tricecane         2.11E-06           2.11E-06         Trimethylbenzene, 12.3-         2.11E-06           2.11E-06         Trimethylbenzene, 12.3-         2.11E-06           2.11E-06         Trimethylbenzene, 13.5-         2.11E-06           2.11E-06         Trimethylbenzene, 13.3-	Biphenyl	2.11E-06	Indene	2.11E-06	Biphe
2.11E-06         Isopropylbenzene         2.11E-06           2.11E-06         Methane         2.30E-04           3.18E-06         Methylnaphthalene, 1-         2.11E-06           6.32E-04         Methylnaphthalene, 2-         2.11E-06           4.02E-04         Norane         2.11E-06           2.11E-06         Octane         2.11E-06           2.11E-06         Octane         2.11E-06           2.11E-06         Phanol         2.11E-06           Porpionaldehyde (Propanal)         1.91E-06           Propylbenzene, n-         2.11E-06           Styrene         2.11E-06           Styrene         2.11E-06           Styrene         2.11E-06           Styrene         2.11E-06           Styrene         2.11E-06           Styrene         2.11E-06           Suffur Dioxide         2.53E-05           Tetadecane         2.11E-06           Tricdecane         2.11E-06           Tricdecane         2.11E-06           Trimethylbenzene, 1,2,3-         2.11E-06           Trimethylbenzene, 1,2,4-         2.11E-06           Trimethylphenol, 2,4,6-         2.11E-06           Trimethylphenol, 2,4,6-         2.11E-06	Butylbenzene, n-	2.11E-06	Isobutylbenzene	2.11E-06	Butyl
2.11E-06       Methane       2.30E-04         3.18E-06       Methylnaphthalene, 1-       2.11E-06         6.32E-04       Methylnaphthalene, 2-       2.11E-06         4.02E-04       Naphthalene       2.11E-06         2.11E-06       Nitrogen Oxides       4.31E-04         2.11E-06       O.m.p-Tolualdehyde       5.09E-06         2.11E-06       Octane       2.11E-06         2.11E-06       Propionaldehyde (Propanal)       1.91E-06         2.11E-06       Propylbenzene, n-       2.11E-06         2.11E-06       Propylbenzene, n-       2.11E-06         2.11E-06       Styrene       2.11E-06         2.11E-06       Styrene       2.11E-06         2.11E-06       Styrene       2.11E-06         2.11E-06       Styrene       2.11E-06         2.11E-06       Tridecane       2.11E-06         2.11E-06       Trimethylbenzene, 1,2,3-       2.11E-06         2.11E-06       Trimethylphenol, 2,4,6-       2.11E-06         2.11E-06       Trimethylphenol, 2,4,6-       2.11E-06         2.11E-06       Trimethylphenol, 2,4,6-       2.11E-06         2.11E-06       Trimethylphenol, 2,4,6-       2.11E-06         2.11E-06       Trimethylphenol, 2,	Butylbenzene, sec-	2.11E-06	Isopropylbenzene	2.11E-06	Buty
3.18E-06       Methylnaphthalene, 1-       2.11E-06         6.32E-04       Methylnaphthalene, 2-       2.11E-06         4.02E-04       Naphthalene       2.11E-06         2.11E-06       Vanane       2.11E-06         2.11E-06       O,m.p-Tolualdehyde       2.11E-06         2.11E-06       Octane       2.11E-06         2.11E-06       Propionaldehyde (Propanal)       1.91E-06         2.11E-06       Propylbenzene, n-       2.11E-06         2.11E-06       Propylbenzene, n-       2.11E-06         2.11E-06       Styrene       2.11E-06         2.11E-06       Styrene       2.11E-06         2.11E-06       Styrene       2.11E-06         2.11E-06       Styrene       2.11E-06         2.11E-06       Tridecane       2.11E-06         2.11E-06       Trimethylbenzene, 1,2,3-       2.11E-06         2.11E-06       Trimethylphenol, 2,3,5-       2.11E-06         2.11E-06       Trimethylphenol, 2,4,6-       2.11E-06         2.11E-06       Trimethylphenol, 2,4,6-       2.11E-06         2.11E-06       Trimethylphenol, 2,4,6-       2.11E-06         2.11E-06       Trimethylphenol, 2,4,6-       2.11E-06         2.11E-06       Trimethylphe	Butylbenzene, tert-	2.11E-06	Methane	2.30E-04	Butyl
6.32E-04       Methylnaphthalene, 2-       2.11E-06         4.02E-04       Naphthalene       2.11E-06         2.11E-06       Nitrogen Oxides       4.31E-04         1.91E-06       O.m.p-Tolualdehyde       2.11E-06         2.11E-06       O.tane       2.11E-06         2.11E-06       Octane       2.11E-06         2.11E-06       Pentanal (Valeraldehyde)       1.91E-06         3-       2.11E-06       Propylbenzene, n-       2.11E-06         3-       2.11E-06       Propylbenzene, n-       2.11E-06         1.3-       2.11E-06       Styrene       2.11E-06         1.5-       2.11E-06       Trimethylbenzene, 1,2,3-       2.11E-06         1.5-       2.11E-06       Trimethylphenol, 2,4,6-       2.11E-06         2.7-       2.11E-06       Trimethylphenol, 2,4,6-       2.11E-06	Butyraldehyde/Methacrolein	3.18E-06		2.11E-06	Butyı
4.02E-04       Naphthalene       2.11E-06         2.11E-06       Nitrogen Oxides       4.31E-04         1.91E-06       O.m.p-Tolualdehyde       2.11E-06         2.11E-06       O.ctane       2.11E-06         2.11E-06       Octane       2.11E-06         2.11E-06       Pentanal (Valeraldehyde)       1.91E-06         2.11E-06       Propionaldehyde (Propanal)       1.91E-06         3-       2.11E-06       Propylbenzene, n-       2.11E-06         1.3-       2.11E-06       Styrene       2.11E-06         1.3-       2.11E-06       Styrene       2.11E-06         1.5-       2.11E-06       Trimethylbenzene, 1,2,3-       2.11E-06         2.3-       2.11E-06       Trimethylphenol, 2,4,6-       2.11E-06         2.7-       2.11E-06       Trimethylphenol, 2,4,6-       2.11E-06         2.11E-06       Trimethylphenol, 2,4,6-       2.11E-06         2.11E-06       Trimethylphenol, 2,4,6-       2.11E-06         2.11E-06       Xylene, np-       2.11E-06         2.11E-06       Trimethylphenol, 2,4,6-       2.11E-06         2.11E-06       Trimethylphenol, 2,4,6-       2.11E-06         2.11E-06       Trimethylphenol, 2,4,6-       2.11E-06      <	Carbon Dioxide	6.32E-04	Methylnaphthalene, 2-	2.11E-06	Carb
2.11E-06       Nitrogen Oxides       4.31E-04         2.11E-06       Nonane       2.11E-06         1.91E-06       0,m,p-Tolualdehyde       5.09E-06         2.11E-06       Octane       2.11E-06         2.11E-06       Pentanal (Valeraldehyde)       1.91E-06         2.11E-06       Propionaldehyde (Propanal)       1.91E-06         3-       2.11E-06       Propylbenzene, n-       2.11E-06         1.3-       2.11E-06       Styrene       2.11E-06         1.3-       2.11E-06       Styrene       2.11E-06         1.5-       2.11E-06       Trimethylbenzene, 1,2,3-       2.11E-06         2.3-       2.11E-06       Trimethylbenzene, 1,3,5-       2.11E-06         2.7-       2.11E-06       Trimethylphenol, 2,4,6-       2.11E-06         2.11E-06       Trimethylphenol, 2,4,6-       2.11E-06         2.11E-06       Trimethylphenol, 2,4,6-       2.11E-06         2.11E-06       Xylene, np-       2.11E-06         2.11E-06       Trimethylphenol, 2,4,6-       2.11E-06         2.11E-06       Trimethylphenol, 2,4,6-       2.11E-06         2.11E-06       Trimethylphenol, 2,4,6-       2.11E-06         2.11E-06       Trimethylphenol, 2,4,6-       2.11E-06	Carbon Monoxide	4.02E-04	Naphthalene	2.11E-06	Carb
2.11E-06       Nonane       2.11E-06         1.91E-06       0,m,p-Tolualdehyde       5.09E-06         2.11E-06       Octane       2.11E-06         2.11E-06       Pentanal (Valeraldehyde)       1.91E-06         2.11E-06       Propionaldehyde (Propanal)       1.91E-06         3-       2.11E-06       Propylbenzene, n-       2.11E-06         1.3-       2.11E-06       Styrene       2.11E-06         1.3-       2.11E-06       Suffur Dioxide       2.11E-06         1.5-       2.11E-06       Trimethylbenzene, 1,2,3-       2.11E-06         1.5-       2.11E-06       Trimethylbenzene, 1,3,5-       2.11E-06         2.7-       2.11E-06       Trimethylbenol, 2,4,6-       2.11E-06         2.7-       2.11E-06       Trimethylbenol, 2,4,6-       2.11E-06         2.11E-06       Trimethylbenol, 2,4,6-       2.11E-06         2.11E-06       Xylene, np-       2.11E-06         2.11E-06       2.11E-06	Cresol, mp-	2.11E-06	Nitrogen Oxides	4.31E-04	Cres
1.91E-06       0,m,p-Tolualdehyde       5.09E-06         2.11E-06       Octane       2.11E-06         2.11E-06       Pentanal (Valeraldehyde)       1.91E-06         2.11E-06       Propionaldehyde (Propanal)       1.91E-06         3.4       2.11E-06       Propylbenzene, n-       2.11E-06         1.3-       2.11E-06       Styrene       2.11E-06         1.3-       2.11E-06       Styrene       2.11E-06         1.5-       2.11E-06       Trimethylbenzene, 1,2,3-       2.11E-06         1.6-       2.11E-06       Trimethylbenzene, 1,2,4-       2.11E-06         2.2-       Trimethylbenzene, 1,3,5-       2.11E-06         2.7-       2.11E-06       Trimethylbenol, 2,3,5-       2.11E-06         2.11E-06       Trimethylbenol, 2,4,6-       2.11E-06         2.11E-06       Trimethylphenol, 2,4,6-       2.11E-06         2.11E-06       Xylene, np-       2.11E-06         2.11E-06       Xylene, np-       2.11E-06	Cresol, o-	2.11E-06	Nonane	2.11E-06	Cres
2.11E-06       Octane       2.11E-06         2.11E-06       Pentanal (Valeraldehyde)       1.91E-06         2.11E-06       Phenol       2.11E-06         2.11E-06       Propionaldehyde (Propanal)       1.91E-06         3-       2.11E-06       2.11E-06         3-       2.11E-06       2.11E-06         1.3-       2.11E-06       2.11E-06         1.5-       2.11E-06       2.11E-06         1.5-       2.11E-06       2.11E-06         1.5-       2.11E-06       2.11E-06         1.5-       2.11E-06       2.11E-06         2.2-1-06       Trimethylbenzene, 1,2,3-       2.11E-06         2.2-1-1-06       Trimethylbenzene, 1,3,5-       2.11E-06         2.7-       2.11E-06       Trimethylbenzene, 1,3,5-       2.11E-06         2.11E-06       Trimethylphenol, 2,4,6-       2.11E-06         2.11E-06       Xylene, mp-       2.11E-06         2.11E-06       2.11E-06	Crotonaldehyde	1.91E-06	o,m,p-Tolualdehyde	5.09E-06	Crotc
2.11E-06       Pentanal (Valeraldehyde)       1.91E-06         2.11E-06       Phenol       2.11E-06         2.11E-06       Propionaldehyde (Propanal)       1.91E-06         3-       2.11E-06       2.11E-06         3-       2.11E-06       2.11E-06         1.3-       2.11E-06       2.11E-06         1.5-       2.11E-06       2.11E-06         1.6-       2.11E-06       2.11E-06         1.6-       2.11E-06       2.11E-06         1.6-       2.11E-06       2.11E-06         2.2-       2.11E-06       2.11E-06         2.11E-06       2.11E-06       2.11E-06         2.11E-06       2.11E-06       2.11E-06         2.11E-06       2.11E-06       2.11E-06         2.11E-06       2.11E-06       2.11E-06	Cyclohexane	2.11E-06	Octane	2.11E-06	Cyclc
2.11E-06       Phenol       2.11E-06         2.11E-06       Propionaldehyde (Propanal)       1.91E-06         3-       2.11E-06       2.11E-06         3-       2.11E-06       2.11E-06         1.3-       2.11E-06       2.11E-06         1.5-       2.11E-06       2.11E-06         2.2-       2.11E-06       2.11E-06         2.7-       2.11E-06       2.11E-06         2.7-       2.11E-06       2.11E-06         2.7-       2.11E-06       2.11E-06	Cymene, p-	2.11E-06	Pentanal (Valeraldehyde)	1.91E-06	Cyme
2.11E-06 Propionaldehyde (Propanal) 1.91E-06 Propionaldehyde (Propanal) 2.11E-06 Styrene 2.11E-06 Styrene 2.11E-06 Styrene 2.11E-06 Suffur Dioxide 2.53E-05 THZ as Propane 2.11E-06 Thradecane 2.11E-06 Trimethylbenzene, 1,2,4 2.11E-06 Trimethylbenzene, 1,2,4 2.11E-06 Trimethylphenol, 2,4,6- 2.11E-0	Decane	2.11E-06	Phenol	2.11E-06	Deca
2.11E-06 Styrene 2.11E-06 Styrene 2.11E-06 Styrene 2.11E-06 Suffur Dioxide 2.53E-05 1.2-11E-06 Suffur Dioxide 2.53E-05 2.11E-06 Thradecane 2.11E-06 Thradecane 2.11E-06 Trinethylbenzene, 1,2,4 2.11E-06 Trimethylbenzene, 1,2,4 2.11E-06 Trimethylbenzene, 1,3,5 2.11E-06 Trimethylbenzene, 1,3,5 2.11E-06 Trimethylbendol, 2,4,6- 2.11E-	Diethylbenzene, 1,2-	2.11E-06	Propionaldehyde (Propanal)	1.91E-06	Dieth
3.4     2.11E-06     Styrene     2.11E-06       1.2-     2.11E-06     Sulfur Dioxide     2.53E-05       1.2-     2.11E-06     Tetradecane     2.11E-06       1.3-     2.11E-06     The as Propane     6.32E-04       1.5-     2.11E-06     Tridecane     2.11E-06       1.6-     2.11E-06     Trimethylbenzene, 1,2,3     2.11E-06       2.3-     2.11E-06     Trimethylbenzene, 1,3,5     2.11E-06       2.7-     2.11E-06     Trimethylphenol, 2,4,6     2.11E-06       2.11E-06     Trimethylphenol, 2,4,6     2.11E-06       2.11E-06     Trimethylphenol, 2,4,6     2.11E-06       2.11E-06     Xylene, mp     2.11E-06       2.11E-06     Xylene, o-     2.11E-06	Diethylbenzene, 1,3-	2.11E-06	Propylbenzene, n-	2.11E-06	Dieth
.3-         2.11E-06         Sulfur Dioxide         2.53E-05           1,2-         2.11E-06         Tetradecane         2.11E-06           1,3-         2.11E-06         THC as Propane         6.32E-04           1,5-         2.11E-06         Toluene         2.11E-06           1,6-         2.11E-06         Trimethylbenzene, 1,2,3         2.11E-06           1,8-         2.11E-06         Trimethylbenzene, 1,3,5         2.11E-06           2,7-         2.11E-06         Trimethylphenol, 2,4,6-         2.11E-06           2,7-         2.11E-06         Trimethylphenol, 2,4,6-         2.11E-06           2,11E-06         Trimethylphenol, 2,4,6-         2.11E-06           2,11E-06         Xylene, mp-         2.11E-06           2,11E-06         Xylene, o-         2.11E-06	Diethylbenzene, 1,4-	2.11E-06	Styrene	2.11E-06	Dieth
1,2-         2.11E-06         Tetradecane         2.11E-06           1,3-         2.11E-06         THC as Propane         6.32E-04           1,5-         2.11E-06         Toluene         2.11E-06           1,6-         2.11E-06         Trimethylbenzene, 1,2,3-         2.11E-06           2,3-         2.11E-06         Trimethylbenzene, 1,2,4-         2.11E-06           2,7-         2.11E-06         Trimethylphenol, 2,4,6-         2.11E-06           2,11E-06         Trimethylphenol, 2,4,6-         2.11E-06           2,11E-06         Trimethylphenol, 2,4,6-         2.11E-06           2,11E-06         Xylene, mp-         2.11E-06           2,11E-06         Xylene, o-         2.11E-06	Diisopropylbenzene, 1,3-	2.11E-06	Sulfur Dioxide	2.53E-05	Diiso
1,3-         2.11E-06         THC as Propane         6.32E-04           1,5-         2.11E-06         Toluene         2.11E-06           1,6-         2.11E-06         Trimethylbenzene, 1,2,3-         2.11E-06           2,3-         2.11E-06         Trimethylbenzene, 1,2,4-         2.11E-06           2,7-         2.11E-06         Trimethylbenzene, 1,3,5-         2.11E-06           2,7-         2.11E-06         Trimethylphenol, 2,4,6-         2.11E-06           2.11E-06         Trimethylphenol, 2,4,6-         2.11E-06           2.11E-06         Xylene, mp-         2.11E-06           2.11E-06         Xylene, o-         2.11E-06		2.11E-06	Tetradecane	2.11E-06	Dime
1,5-         2.11E-06         Toluene         2.11E-06           1,6-         2.11E-06         Tridecane         2.11E-06           1,8-         2.11E-06         Trimethylbenzene, 1,2,3         2.11E-06           2,3-         2.11E-06         Trimethylbenzene, 1,3,5         2.11E-06           2,7-         2.11E-06         Trimethylphenol, 2,3,5-         2.11E-06           2.11E-06         Trimethylphenol, 2,4,6-         2.11E-06           2.11E-06         Trimethylphenol, 2,4,6-         2.11E-06           2.11E-06         Xylene, mp-         2.11E-06           2.11E-06         Xylene, o-         2.11E-06           2.11E-06         2.11E-06		2.11E-06	THC as Propane	6.32E-04	Dime
1,6-       2.11E-06       Tridecane       2.11E-06         1,8-       2.11E-06       Trimethylbenzene, 1,2,3-       2.11E-06         2,3-       2.11E-06       Trimethylbenzene, 1,2,4-       2.11E-06         2,6-       2.11E-06       Trimethylbenzene, 1,3,5-       2.11E-06         2,7-       2.11E-06       Trimethylphenol, 2,3,5-       2.11E-06         2.11E-06       Trimethylphenol, 2,4,6-       2.11E-06         2.11E-06       Xylene, mp-       2.11E-06         2.11E-06       Xylene, o-       2.11E-06	Dimethylnaphthalene, 1,5-	2.11E-06	Toluene	2.11E-06	Dime
1,8-     2.11E-06     Trimethylbenzene, 1,2,3-     2.11E-06       2,3-     2.11E-06     Trimethylbenzene, 1,2,4-     2.11E-06       2,6-     2.11E-06     Trimethylbenzene, 1,3,5-     2.11E-06       2,7-     2.11E-06     Trimethylphenol, 2,3,5-     2.11E-06       2.11E-06     Trimethylphenol, 2,4,6-     2.11E-06       2.11E-06     Wylene, mp-     2.11E-06       2.11E-06     Xylene, o-     2.11E-06	Dimethylnaphthalene, 1,6-	2.11E-06	Tridecane	2.11E-06	Dime
2,3-       2.11E-06       Trimethylbenzene, 1,2,4-       2.11E-06         2,6-       2.11E-06       Trimethylbenzene, 1,3,5-       2.11E-06         2,7-       2.11E-06       Trimethylphenol, 2,3,5-       2.11E-06         2.11E-06       Trimethylphenol, 2,4,6-       2.11E-06         2.11E-06       Lndecane       2.11E-06         2.11E-06       Xylene, mp-       2.11E-06         2.11E-06       Xylene, o-       2.11E-06	Dimethylnaphthalene, 1,8-	2.11E-06	Trimethylbenzene, 1,2,3-	2.11E-06	Dime
2,6-       2.11E-06       Trimethylbenzene, 1,3,5-       2.11E-06         2,7-       2.11E-06       Trimethylphenol, 2,3,5-       2.11E-06         2.11E-06       Trimethylphenol, 2,4,6-       2.11E-06         2.11E-06       Wylene, mp-       2.11E-06         2.11E-06       Xylene, o-       2.11E-06         2.11E-06       Xylene, o-       2.11E-06	Dimethylnaphthalene, 2,3-	2.11E-06	Trimethylbenzene, 1,2,4-	2.11E-06	Dime
2,7-     2.11E-06     Trimethylnaphthalene, 2,3,5-     2.11E-06       2.11E-06     Trimethylphenol, 2,3,6-     2.11E-06       2.11E-06     Undecane     2.11E-06       2.11E-06     Xylene, o-     2.11E-06       2.11E-06     Xylene, o-     2.11E-06	Dimethylnaphthalene, 2,6-	2.11E-06	Trimethylbenzene, 1,3,5-	2.11E-06	Dime
2.11E-06 Trimethylphenol, 2,3,5- 2.11E-06 2.11E-06 Trimethylphenol, 2,4,6- 2.11E-06 2.11E-06 Xylene, mp- 2.11E-06 Xylene, o- 2.11E-06 2.11E-06 Xylene, o- 2.11E-06 2.11E-06 Xylene, o- 2.11E-06	Dimethylnaphthalene, 2,7-	2.11E-06	Trimethylnaphthalene, 2,3,5-	2.11E-06	Dime
2.11E-06 Trimethylphenol, 2,4,6- 2.11E-06 Undecane 2.11E-06 Xylene, mp- 2.11E-06 Xylene, o- 2.11E-06 Xylen	Dimethylphenol, 2,3-	2.11E-06	Trimethylphenol, 2,3,5-	2.11E-06	Dime
2.11E-06     Undecane     2.11E-06       2.11E-06     Xylene, mp-     2.11E-06       2.11E-06     Xylene, o-     2.11E-06	Dimethylphenol, 2,4-	2.11E-06	Trimethylphenol, 2,4,6-	2.11E-06	Dime
2.11E-06 Xylene, mp- 2.11E-06 Xylene, o- 2.11E-06 Zylene, o- 2.11E	Dimethylphenol, 2,5-	2.11E-06	Undecane	2.11E-06	Dime
2.11E-06 Xylene, o- 2.11E-06 2.11E-06 2.11E-06	Dimethylphenol, 2,6-	2.11E-06	Xylene, mp-	2.11E-06	DIMe
2.11E-06	Dimethylphenol, 3,4-	2.11E-06	Xylene, o-	2.11E-06	e   c
	Dimethylphenol, 3,5-	2.11E-06			Dime

Analyte	lb/tn core	Analyte	lb/tn core
2-Butanone (MEK)	1.11E-04	Dodecane	1.23E-04
Acenaphthalene	1.23E-04	Ethylbenzene	1.23E-04
Acetaldehyde	1.11E-04	Ethyltoluene, 2-	1.23E-04
Acetophenone	1.43E-02	Ethyltoluene, 3-	1.23E-04
Acrolein	1.11E-04	Ethyltoluene, 4-	1.23E-04
alpha-Methylstyrene	1.23E-04	Formaldehyde	1.11E-04
Ammonia	4.32E-03	Heptane	1.23E-04
Anthracene	1.23E-04	Hexaldehyde	1.11E-04
Benzaldehyde	1.11E-04	Hexane	1.23E-04
Benzene	1.23E-04	Indan	1.23E-04
Biphenyl	1.23E-04	Indene	1.23E-04
Butylbenzene, n-	1.23E-04	Isobutylbenzene	1.23E-04
Butylbenzene, sec-	1.23E-04	Isopropylbenzene	1.23E-04
Butylbenzene, tert-	1.23E-04	Methane	1.34E-02
Butyraldehyde/Methacrolein	1.86E-04	Methylnaphthalene, 1-	1.23E-04
Carbon Dioxide	3.68E-02	Methylnaphthalene, 2-	1.23E-04
Carbon Monoxide	2.34E-02	Naphthalene	1.23E-04
Cresol, mp-	1.23E-04	Nitrogen Oxides	2.51E-02
Cresol, o-	1.23E-04	Nonane	1.23E-04
Crotonaldehyde	1.11E-04	o,m,p-Tolualdehyde	2.97E-04
Cyclohexane	1.23E-04	Octane	1.23E-04
Cymene, p-	1.23E-04	Pentanal (Valeraldehyde)	1.11E-04
Decane	1.23E-04	Phenol	1.23E-04
	1.23E-04	Propionaldehyde (Propanal)	1.11E-04
Diethylbenzene, 1,3-	1.23E-04	Propylbenzene, n-	1.23E-04
Diethylbenzene, 1,4-	1.23E-04	Styrene	1.23E-04
	1.23E-04	Sulfur Dioxide	1.48E-03
Dimethylnaphthalene, 1,2-	1.23E-04	Tetradecane	1.23E-04
	1.23E-04	THC as Propane	3.68E-02
	1.23E-04	Toluene	1.23E-04
	1.23E-04	Tridecane	1.23E-04
	1.23E-04	Trimethylbenzene, 1,2,3-	1.23E-04
Dimethylnaphthalene, 2,3-	1.23E-04	Trimethylbenzene, 1,2,4-	1.23E-04
	1.23E-04	꺆	1.23E-04
Dimethylnaphthalene, 2,7-	1.23E-04	Trimethylnaphthalene, 2,3,5-	1.23E-04
Dimethylphenol, 2,3-	1.23E-04	Trimethylphenol, 2,3,5-	1.23E-04
Dimethylphenol, 2,4-	1.23E-04	Trimethylphenol, 2,4,6-	1.23E-04
Dimethylphenol, 2,5-	1.23E-04	Undecane	1.23E-04
Dimethylphenol, 2,6-	1.23E-04	Xylene, mp-	1.23E-04
Dimethylphenol, 3,4-	1.23E-04	Xylene, o-	1.23E-04
Dimethylphenol, 3,5-	1.23E-04		

Detailed Emissions - Core Make/Store and Storage, Test HO, Ib/ton core

			Detailed Emiss	ions - Co	re Make/S	Store and	Storage,	Test HO,	lb/ton co	re	
TA	POM	НАР		HO001	HO002	HO003	HO004	HO005	HO006	Average	Standard Deviation
_		-	Test Dates	25-Apr-07	25-Apr-07	26-Apr-07	26-Apr-07	26-Apr-07	26-Apr-07	-	_
Emi	ssio	n Ind	icators								
			THC as Propane	8.21E-01	9.94E-01	1.09E+00	8.17E-01	8.12E-01	1.08E+00	9.34E-01	1.33E-01
			Non-Methane Hydrocarbons	4.42E-01	4.40E-01	1.82E-01	1.40E-01	1.52E-01	4.21E-01	2.96E-01	1.52E-01
			Sum of Target Analytes	1.35E-01	1.44E-01	1.53E-01	1.53E-01	1.45E-01	1.45E-01	1.46E-01	6.86E-03
			Sum of Target HAPs	3.90E-02	3.93E-02	3.93E-02	4.50E-02	4.11E-02	4.01E-02	4.06E-02	2.27E-03
			Sum of Target POMs	4.99E-05	6.00E-05	4.25E-05	3.73E-05	4.94E-05	4.25E-05	4.69E-05	7.97E-06
Sele	cted	Tar	get HAPs and POMs		****		****				
TA			Phenol	3.15E-02	3.13E-02	3.08E-02	3.48E-02	3.34E-02	3.23E-02	3.23E-02	1.51E-03
TA		Н	Formaldehyde	6.60E-03	6.77E-03	7.40E-03	8.97E-03	6.47E-03	6.88E-03	7.18E-03	9.35E-04
TA		Н	Styrene	1.89E-04	4.21E-04	2.83E-04	2.44E-04	4.60E-04	2.12E-04	3.01E-04	1.13E-04
TA			Acetaldehyde	2.19E-04	2.62E-04	2.25E-04	3.80E-04	2.54E-04	2.33E-04	2.62E-04	6.00E-05
TA		Н	Propionaldehyde (Propanal)	1.67E-04	1.90E-04	1.97E-04	1.86E-04	1.60E-04	1.78E-04	1.79E-04	1.40E-05
TA		_	Acrolein	1.25E-04	1.62E-04	1.59E-04	1.67E-04	1.74E-04	1.36E-04	1.54E-04	1.92E-05
TA		Н	Hexane	5.87E-05	5.14E-05	8.79E-05	9.40E-05	≤PQL	6.25E-05	6.38E-05	2.43E-05
TA	Р	Н	Naphthalene	4.99E-05	6.00E-05	4.25E-05	3.73E-05	4.94E-05	4.25E-05	4.69E-05	7.97E-06
TA		Н	Toluene	3.73E-05	≤PQL	3.42E-05	4.04E-05	≤PQL	≤PQL	3.27E-05	5.36E-06
TA		Н	Cresol, o-	4.87E-05	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	3.16E-05	8.39E-06
TA		Н	Benzene	≤PQL	≤PQL	≤PQL	3.01E-05	3.17E-05	3.07E-05	2.95E-05	1.54E-06
TA	Р	Н	Acenaphthalene	≤PQL       NA							
TA	Р	Н	Anthracene	≤PQL       NA							
TA	P	Н	Dimethylnaphthalene, 1,2-	≤PQL       NA							
TA	P	Н	Dimethylnaphthalene, 1,3-	≤PQL       NA							
TA	Р	Н	Dimethylnaphthalene, 1,5-	≤PQL       NA							
TA	Р	Н	Dimethylnaphthalene, 1,6-	≤PQL       NA							
TA	Р	Н	Dimethylnaphthalene, 1,8-	≤PQL       NA							
TA	Р	Н	Dimethylnaphthalene, 2,3-	≤PQL       NA							
TA	Р	н	Dimethylnaphthalene, 2,6-	≤PQL       NA							
TA	Р	Н	Dimethylnaphthalene, 2,7-	≤PQL       NA							
TA	Р	Н	Methylnaphthalene, 1-	≤PQL       NA							
TA	Р	Н	Methylnaphthalene, 2-	≤PQL       NA							
TA	Р	Н	Trimethylnaphthalene, 2,3,5-	≤PQL       NA							
TA		Н	Biphenyl	≤PQL       NA							
TA		Н	Cresol, mp-	ı	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA		Н	Ethylbenzene	≤PQL       NA							
TA		Н	Xylene, mp-	≤PQL       NA							
TA		Н	Xylene, o-	≤PQL       NA							
TA			Acetophenone	≤PQL       NA							
Add	ition	al Se	elected Target Analytes				-				
TA			Ammonia	9.40E-02	1.03E-01	1.12E-01	1.07E-01	1.02E-01	1.03E-01	1.03E-01	5.95E-03
TA			Crotonaldehyde	7.95E-04	8.93E-04	1.02E-03	8.87E-04	8.43E-04	7.84E-04	8.70E-04	8.50E-05
TA			2-Butanone (MEK)	2.89E-04	3.19E-04	3.53E-04	2.43E-04	3.11E-04	2.93E-04	3.02E-04	3.67E-05
TA			Butyraldehyde/Methacrolein	2.96E-04	3.43E-04	3.11E-04	2.47E-04	3.26E-04	2.41E-04	2.94E-04	4.18E-05
TA			Dimethylphenol, 3,4-	3.40E-04	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	8.01E-05	1.27E-04
TA			Trimethylbenzene, 1,2,4-	4.08E-05	4.02E-05	5.66E-05	5.09E-05	4.34E-05	4.43E-05	4.60E-05	6.45E-06
TA			Heptane	3.35E-05	3.54E-05	4.45E-05	5.77E-05	3.92E-05	4.24E-05	4.21E-05	8.68E-06
TA			Diethylbenzene, 1,4-	8.38E-05	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	3.74E-05	2.27E-05
TA			Diethylbenzene, 1,3-	4.07E-05	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	3.02E-05	5.12E-06
TA			Butylbenzene, sec-	3.87E-05	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	2.99E-05	4.30E-06

(continued on next page)

ND=Not Detected NA= Not Applicable I=Invalidated Data

≤PQL=Less than or equal to the Practical Quantitation Limit

Detailed Emissions - Core Make/Store and Storage, Test HO, lb/ton core (continued)

								1001110,		(0011011	Standard
ΔT	POM	ΗΨ		HO001	HO002	HO003	HO004	HO005	HO006	Average	Deviation
_		<del>  _</del>	Test Dates	25-Apr-07	25-Apr-07	26-Apr-07	26-Apr-07	26-Apr-07	26-Apr-07	- Average	-
TA			Tetradecane	≤PQL	3.02E-05	≤PQL	≤PQL	2.96E-05	≤PQL	2.87E-05	9.02E-07
TA			Benzaldehyde	≤PQL	1	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA			Cyclohexane	≤PQL	NA						
TA			Decane	≤PQL	NA						
TA			Dimethylphenol, 2,4-	≤PQL	NA						
TA			Dimethylphenol, 2,6-	≤PQL	NA						
TA			Dodecane	≤PQL	NA						
TA			Ethyltoluene, 2-	1	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA			Ethyltoluene, 3-	≤PQL	NA						
TA			Hexaldehyde	≤PQL	NA						
TA			Indan	≤PQL	NA						
TA			Nonane	1	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA			o,m,p-Tolualdehyde	≤PQL	NA						
TA			Octane	≤PQL	NA						
TA			Pentanal (Valeraldehyde)	≤PQL	NA						
TA			Propylbenzene, n-	≤PQL	NA						
TA			Trimethylbenzene, 1,2,3-	≤PQL	NA						
TA			Trimethylbenzene, 1,3,5-	≤PQL	NA						
TA			Undecane	≤PQL	NA						
TA			alpha-Methylstyrene	≤PQL	NA						
TA			Butylbenzene, n-	≤PQL	NA						
TA			Butylbenzene, tert-	≤PQL	NA						
TA			Cymene, p-	≤PQL	NA						
TA			Diethylbenzene, 1,2-	≤PQL	NA						
TA			Diisopropylbenzene, 1,3-	≤PQL	NA						
TA			Dimethylphenol, 2,3-	≤PQL	NA						
TA			Dimethylphenol, 2,5-	≤PQL	NA						
TA			Dimethylphenol, 3,5-	≤PQL	NA						
TA			Ethyltoluene, 4-	≤PQL	NA						
TA			Indene	≤PQL	NA						
TA			Isobutylbenzene	≤PQL	NA						
TA			Isopropylbenzene	≤PQL	NA						
TA			Tridecane	≤PQL	NA						
TA			Trimethylphenol, 2,3,5-	≤PQL	NA						
TA			Trimethylphenol, 2,4,6-	≤PQL	NA						
Sele	cted	Crit	eria Pollutants and Greenhous								
			Carbon Dioxide	1.42E+01	1.58E+01	1.69E+01	1.44E+01	1.41E+01	1.40E+01	1.49E+01	1.19E+00
			Methane	3.79E-01	5.53E-01	9.04E-01	6.77E-01	6.59E-01	6.57E-01	6.38E-01	1.72E-01
			Carbon Monoxide	1.15E-01	9.87E-02	1.35E-01	1.34E-01	1.05E-01	1.02E-01	1.15E-01	1.62E-02
			Nitrogen Oxides	≤PQL	NA						
			Sulfur Dioxide	≤PQL	NA						

ND=Not Detected NA= Not Applicable I=Invalidated Data

≤PQL=Less than or equal to the Practical Quantitation Limit

## Detailed Emissions Core Make/Store and Storage, Test HO, lb/lb binder

			Detailed Eilligg				- · · · · · · · · · · · · · · · · · · ·				
₹	POM	₽		HO001	HO002	HO003	HO004	HO005	HO006	Average	Standard Deviation
	_	_	Test Dates	25-Apr-07	25-Apr-07	26-Apr-07	26-Apr-07	26-Apr-07	26-Apr-07	_	_
Fmi	ssio	n Ind	icators	20710.01	20740.0.	20740.01	20740.0.	20710.0.	20740. 01		
		l	THC as Propane	1.41E-02	1.70E-02	1.87E-02	1.40E-02	1.39E-02	1.85E-02	1.60E-02	2.29E-03
			Non-Methane Hydrocarbons	7.58E-03	7.56E-03	3.12E-03	2.40E-03	2.61E-03	7.22E-03	5.08E-03	2.61E-03
			Sum of Target Analytes	2.32E-03	2.47E-03	2.62E-03	2.71E-03	2.54E-03	2.54E-03	2.53E-03	1.35E-04
			Sum of Target HAPs	6.71E-04	6.74E-04	6.72E-04	7.96E-04	7.22E-04	7.05E-04	7.07E-04	4.85E-05
			Sum of Target POMs	8.55E-07	1.03E-06	7.28E-07	6.61E-07	8.67E-07	7.46E-07	8.15E-07	1.32E-07
Sele	cted	Tar	get HAPs and POMs	0.002 0.	1.002 00		0.0.2	0.0.2 0.		0.102 01	
TA			Phenol	5.39E-04	5.37E-04	5.27E-04	6.16E-04	5.87E-04	5.67E-04	5.62E-04	3.45E-05
TA		_	Formaldehyde	1.16E-04	1.16E-04	1.27E-04	1.59E-04	1.14E-04	1.21E-04	1.25E-04	1.71E-05
TA		Н	Styrene	3.23E-06	7.23E-06	4.85E-06	4.32E-06	8.08E-06	3.72E-06	5.24E-06	1.97E-06
TA		Н	Acetaldehyde	3.83E-06	4.49E-06	3.85E-06	6.73E-06	4.46E-06	4.09E-06	4.58E-06	1.09E-06
TA		Н	Propionaldehyde (Propanal)	3.17E-06	3.25E-06	3.37E-06	3.29E-06	2.81E-06	3.12E-06	3.17E-06	1.96E-07
TA		H	Acrolein	2.42E-06	2.77E-06	2.73E-06	2.97E-06	3.05E-06	2.38E-06	2.72E-06	2.74E-07
TA		_	Hexane	1.01E-06	8.82E-07	1.50E-06	1.67E-06	≤PQL	1.10E-06	1.11E-06	4.29E-07
TA	Р	Н	Naphthalene	8.55E-07	1.03E-06	7.28E-07	6.61E-07	8.67E-07	7.46E-07	8.15E-07	1.32E-07
TA	•	H	Toluene	6.38E-07	≤PQL	5.85E-07	7.15E-07	≤PQL	≤PQL	5.64E-07	9.84E-08
TA		H	Cresol, o-	8.35E-07	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	5.42E-07	1.44E-07
TA		_	Benzene	≤PQL	≤PQL	≤PQL	5.33E-07	5.56E-07	5.39E-07	5.13E-07	3.36E-08
TA	Р	_	Acenaphthalene	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA	P	Н	Anthracene	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA NA
TA	P		Dimethylnaphthalene, 1,2-	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA NA
TA	P		Dimethylnaphthalene, 1,3-	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA NA
TA	P	_	Dimethylnaphthalene, 1,5-	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA NA
TA	P		Dimethylnaphthalene, 1,6-	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA NA
TA	P	H	Dimethylnaphthalene, 1,8-	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA NA
TA	P	_	Dimethylnaphthalene, 2,3-	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA NA
TA	P		Dimethylnaphthalene, 2,6-	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA NA
TA	P		Dimethylnaphthalene, 2,7-	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA NA
TA	P		Methylnaphthalene, 1-	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA NA
TA	P		Methylnaphthalene, 2-	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA NA
TA	P	H	Trimethylnaphthalene, 2,3,5-	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA NA
TA	-	H	Biphenyl	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA NA
TA		H	Cresol, mp-	1	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA NA
TA		H	Ethylbenzene	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA NA
TA		_	Xylene, mp-	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA NA
TA		H	Xylene, o-	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA NA
TA			Acetophenone	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
	ition	al Se	elected Target Analytes	=1 &L	i &L		QL	i &L	_1 QL	_1 QL	14/1
TA		<u> </u>	Ammonia	1.61E-03	1.77E-03	1.92E-03	1.89E-03	1.79E-03	1.81E-03	1.80E-03	1.08E-04
TA			Crotonaldehyde	1.48E-05	1.53E-05	1.74E-05	1.57E-05	1.48E-05	1.38E-05	1.53E-05	1.22E-06
TA			Butyraldehyde/Methacrolein	5.87E-06	5.89E-06	5.33E-06	4.37E-06	5.73E-06	4.24E-06	5.24E-06	7.51E-07
TA			2-Butanone (MEK)	4.96E-06	5.48E-06	6.05E-06	4.31E-06	5.46E-06	5.14E-06	5.23E-06	5.87E-07
TA			Dimethylphenol, 3,4-	5.82E-06	5.40L-00 ≤PQL	0.03L-00 ≤PQL	4.51L-00 ≤PQL	5.40L-00 ≤PQL	5.14L-00 ≤PQL	1.37E-06	2.18E-06
TA			Trimethylbenzene, 1,2,4-	7.00E-07	6.89E-07	9.70E-07	9.02E-07	7.62E-07	7.78E-07	8.00E-07	1.13E-07
TA			Heptane	5.75E-07	6.08E-07	7.62E-07	1.02E-06	6.88E-07	7.44E-07	7.33E-07	1.60E-07
TA			Diethylbenzene, 1,4-	1.44E-06	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	6.42E-07	3.89E-07
TA			Diethylbenzene, 1,3-	6.97E-07	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	5.19E-07	8.76E-08
TA			Butylbenzene, sec-	6.63E-07	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	5.13E-07	7.35E-08
TA			Tetradecane	≤PQL	5.17E-07	≤PQL	≤PQL	5.19E-07	≤PQL	4.95E-07	1.83E-08
17A			Totadoodilo		J.17 L 07			J. 10L 01		ed on nex	

(continued on next page)

ND=Not Detected NA= Not Applicable I=Invalidated Data

≤PQL=Less than or equal to the Practical Quantitation Limit

## Detailed Emissions Core Make/Store and Storage, Test HO, lb/lb binder (continued)

			Detailed Lilliss	10113 001	- Water	lore aria t	Jiorage,	1031110,	D/ID DIIIG	Ci (OOiitii	
Ι¥	POM	HAP		HO001	HO002	HO003	HO004	HO005	HO006	Average	Standard Deviation
			Test Dates	25-Apr-07	25-Apr-07	26-Apr-07	26-Apr-07	26-Apr-07	26-Apr-07	- "	_
TA			Cymene, p-	≤PQL	≤PQL	≤PQL	≤PQL	4.90E-07	≤PQL	4.84E-07	2.89E-09
TA			Benzaldehyde	≤PQL		≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA			Cyclohexane	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA			Decane	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA			Dimethylphenol, 2,4-	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA			Dimethylphenol, 2,6-	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA			Dodecane	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA			Ethyltoluene, 2-		≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA			Ethyltoluene, 3-	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA			Hexaldehyde	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA			Indan	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA			Nonane	I	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA			o,m,p-Tolualdehyde	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA			Octane	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA			Pentanal (Valeraldehyde)	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA			Propylbenzene, n-	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA			Trimethylbenzene, 1,2,3-	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA			Trimethylbenzene, 1,3,5-	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA			Undecane	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA			alpha-Methylstyrene	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA			Butylbenzene, n-	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA			Butylbenzene, tert-	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA			Diethylbenzene, 1,2-	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA			Diisopropylbenzene, 1,3-	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA			Dimethylphenol, 2,3-	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA			Dimethylphenol, 2,5-	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA			Dimethylphenol, 3,5-	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA			Ethyltoluene, 4-	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA			Indene	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA			Isobutylbenzene	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA			Isopropylbenzene	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA			Tridecane	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA			Trimethylphenol, 2,3,5-	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
TA			Trimethylphenol, 2,4,6-	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
Sele	cted	Crit	eria Pollutants and Greenhous								
			Carbon Dioxide	2.43E-01	2.71E-01	2.91E-01	2.47E-01	2.42E-01	2.39E-01	2.56E-01	2.06E-02
			Methane	6.50E-03	9.49E-03	1.55E-02	1.16E-02	1.13E-02	1.13E-02	1.10E-02	2.96E-03
			Carbon Monoxide	1.97E-03	1.69E-03	2.32E-03	2.31E-03	1.80E-03	1.75E-03	1.97E-03	2.78E-04
			Nitrogen Oxides	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA
			Sulfur Dioxide	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	≤PQL	NA

ND=Not Detected NA= Not Applicable I=Invalidated Data

 $\leq$ PQL=Less than or equal to the Practical Quantitation Limit

Practical Reporting Limits, Core Make/Store and Storage, Test HO, Ib/Ib binder

Analyte

2-Butanone (MEK)

cenaphthalene

Acetophenone

Ipha-Methylstyrene

Acrolein

Practical Reporting Limits, Core Make/Store and Storage, Test HO, Ib/ton core

lb/lb Binder	Analyte	lb/lb Binder	Analyte	lb/tn core	Analyte	lb/tn core
1.81E-06	Dodecane	4.83E-07	2-Butanone (MEK)	1.05E-04	Dodecane	2.82E-05
4.83E-07	Ethylbenzene	4.83E-07	Acenaphthalene	2.82E-05	Ethylbenzene	2.82E-05
1.81E-06	Ethyltoluene, 2-	4.83E-07	Acetaldehyde	1.05E-04	Ethyltoluene, 2-	2.82E-05
5.88E-05	Ethyltoluene, 3-	4.83E-07	Acetophenone	3.43E-03	Ethyltoluene, 3-	2.82E-05
1.81E-06	Ethyltoluene, 4-	4.83E-07	Acrolein	1.05E-04	Ethyltoluene, 4-	2.82E-05
4.83E-07	Formaldehyde	1.81E-06	alpha-Methylstyrene	2.82E-05	Formaldehyde	1.05E-04
6.07E-05	Heptane	4.83E-07	Ammonia	7.08E-03	Heptane	2.82E-05
3E-07	Hexaldehyde	1.81E-06	Anthracene	2.82E-05	Hexaldehyde	1.05E-04
1.81E-06	Hexane	4.83E-07	Benzaldehyde	1.05E-04	Hexane	2.82E-05
4.83E-07	Indan	4.83E-07	Benzene	2.82E-05	Indan	2.82E-05
E-07	Indene	4.83E-07	Biphenyl	2.82E-05	Indene	2.82E-05
4.83E-07	Isobutylbenzene	4.83E-07	Butylbenzene, n-	2.82E-05	Isobutylbenzene	2.82E-05
E-07	Isopropylbenzene	4.83E-07	Butylbenzene, sec-	2.82E-05	Isopropylbenzene	2.82E-05
4.83E-07	Methane	2.18E-04	Butylbenzene, tert-	2.82E-05	Methane	2.18E-04
3.01E-06	Methylnaphthalene, 1-	4.83E-07	Butyraldehyde/Methacrolein	1.75E-04	Methylnaphthalene, 1-	2.82E-05
5.99E-04	Methylnaphthalene, 2-	4.83E-07	Carbon Dioxide	5.99E-04	Methylnaphthalene, 2-	2.82E-05
3.81E-04	Naphthalene	4.83E-07	Carbon Monoxide	3.81E-04	Naphthalene	2.82E-05
4.83E-07	Nitrogen Oxides	4.08E-04	Cresol, mp-	2.82E-05	Nifrogen Oxides	4.08E-04
E-07	Nonane	4.83E-07	Cresol, o-	2.82E-05	Nonane	2.82E-05
1.81E-06	o,m,p-Tolualdehyde	4.82E-06	Crotonaldehyde	1.05E-04	o,m,p-Tolualdehyde	2.81E-04
4.83E-07	Octane	4.83E-07	Cyclohexane	2.82E-05	Octane	2.82E-05
4.83E-07	Pentanal (Valeraldehyde)	1.81E-06	Cymene, p-	2.82E-05	Pentanal (Valeraldehyde)	1.05E-04
4.83E-07	Phenol	4.83E-07	Decane	2.82E-05	Phenol	2.82E-05
4.83E-07	Propionaldehyde (Propanal)	1.81E-06	Diethylbenzene, 1,2-	2.82E-05	Propionaldehyde (Propanal)	1.05E-04
E-07	Propylbenzene, n-	4.83E-07	Diethylbenzene, 1,3-	2.82E-05	Propylbenzene, n-	2.82E-05
4.83E-07	Styrene	4.83E-07	Diethylbenzene, 1,4-	2.82E-05	Styrene	2.82E-05
4.83E-07	Sulfur Dioxide	2.43E-05	Diisopropylbenzene, 1,3-	2.82E-05	Sulfur Dioxide	1.41E-03
4.83E-07	Tetradecane	4.83E-07	Dimethylnaphthalene, 1,2-	2.82E-05	Tetradecane	2.82E-05
4.83E-07	THC as Propane	5.99E-04	Dimethylnaphthalene, 1,3-	2.82E-05	THC as Propane	6.32E-04
4.83E-07	Toluene	4.83E-07	Dimethylnaphthalene, 1,5-	2.82E-05	Toluene	2.82E-05
4.83E-07	Tridecane	4.83E-07	Dimethylnaphthalene, 1,6-	2.82E-05	Tridecane	2.82E-05
4.83E-07	Trimethylbenzene, 1,2,3-	4.83E-07	Dimethylnaphthalene, 1,8-	2.82E-05	Trimethylbenzene, 1,2,3-	2.82E-05
4.83E-07	Irimethylbenzene, 1,2,4-	4.83E-07	Dimethylnaphthalene, 2,3-	2.82E-05	Trimethylbenzene, 1,2,4-	2.82E-05
4.83E-07	Irimethylbenzene, 1,3,5-	4.83E-07	Dimethylnaphthalene, 2,6-	2.82E-05	Trimethylbenzene, 1,3,5-	2.82E-05
4.83E-07	Irimethylnaphthalene, 2,3,5-	4.83E-07	Dimethylnaphthalene, 2,7-	2.82E-05	Trimethylnaphthalene, 2,3,5-	2.82E-05
4.83E-07	Trimethylphenol, 2,3,5-	4.83E-07	Dimethylphenol, 2,3-	2.82E-05	Trimethylphenol, 2,3,5-	2.82E-05
4.83E-0/	I rimethylphenol, 2,4,6-	4.83E-07	Dimethylphenol, 2,4-	2.82E-05	Trimethylphenol, 2,4,6-	2.82E-05
4.83E-07	Undecane	4.83E-07	Dimethylphenol, 2,5-	2.82E-05	Undecane	2.82E-05
4.83E-07	Xylene, mp-	4.83E-07	Dimethylphenol, 2,6-	2.82E-05	Xylene, mp-	2.82E-05
4.83E-07	Xylene, o-	4.83E-07	Dimethylphenol, 3,4-	2.82E-05	Xylene, o-	2.82E-05
4.83E-0/			i			

Crotonaldehyde Cyclohexane

Jymene, p-

Decane

Dimethylnaphthalene, 1,5-Dimethylnaphthalene, 1,6-

Jimethylnaphthalene, 1,3-

Dimethylnaphthalene, 1,8-Dimethylnaphthalene, 2,3Dimethylnaphthalene, 2,6-Dimethylnaphthalene, 2,7-

Dimethylphenol, 2,3Dimethylphenol, 2,4Dimethylphenol, 2,5Dimethylphenol, 2,6Dimethylphenol, 3,4Dimethylphenol, 3,4-

Dimethylnaphthalene, 1,2-

Diisopropylbenzene, 1,3-

Diethylbenzene, 1,3-Diethylbenzene, 1,4-

Diethylbenzene, 1,2-

Butyraldehyde/Methacrolein

Sarbon Monoxide

Cresol, mp-

Cresol, o-

Carbon Dioxide

Butylbenzene, sec-Butylbenzene, tert-

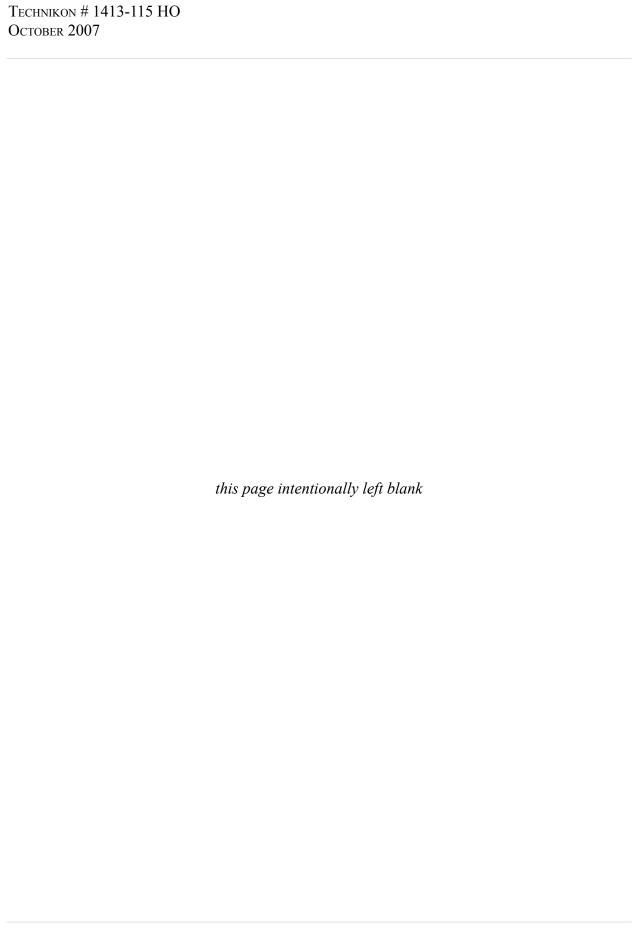
Butylbenzene, n-

Benzaldehyde

Benzene

Bipheny

Ammonia Anthracene APPENDIX C DETAILED PROCESS DATA

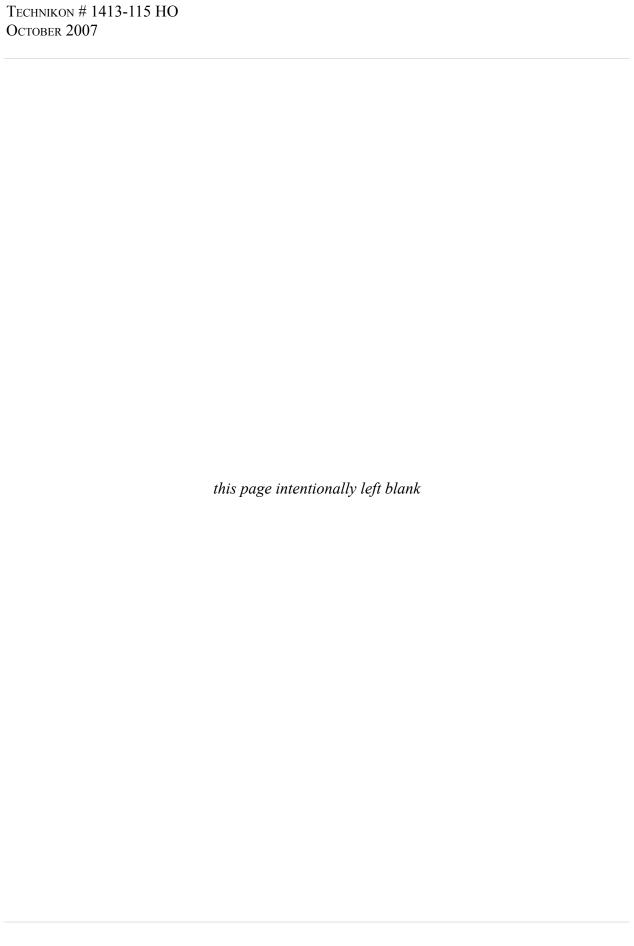


## **Detailed Process Data - Test Series HN**

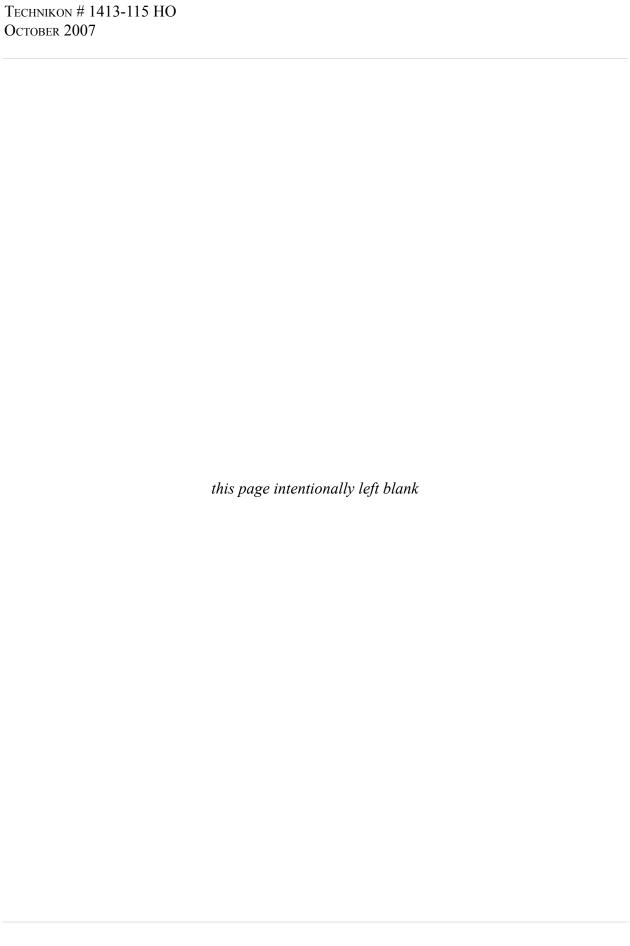
	Detailet	1110003	Data - I	cot ocnic	, , , , ,		
	Tech	nisand XC30	Coremaking				
Test Dates	04/23/07	04/23/07	04/24/07	04/24/07	04/24/07	04/25/07	
Emissions Sample #	HN001	HN002	HN003	HN004	HN005	HN006	Averages
Production Sample #	HN001	HN002	HN003	HN004	HN005	HN006	
Average cooled core weight, g	2479.8	2533.7	2508.8	2539.2	2545.6	2486.5	2515.6
Average core box temperature	452	452	459	460	458	455	456
% stated core binder (BOS)	3	3	3	3	3	3	3
% calculated actual binder	2.9	2.9	2.9	2.9	2.9	2.9	2.9
Number of cores made in run	15	15	15	15	15	15	15
Total core weight in from run, Lbs.	82.01	83.79	82.96	83.97	84.18	82.23	83.19
Total binder weight in run, Lbs.	2.4	2.4	2.4	2.4	2.5	2.4	2.4
Core LOI, %	3.1627	3.2267	3.2703	3.2101	3.2257	3.1912	3.21

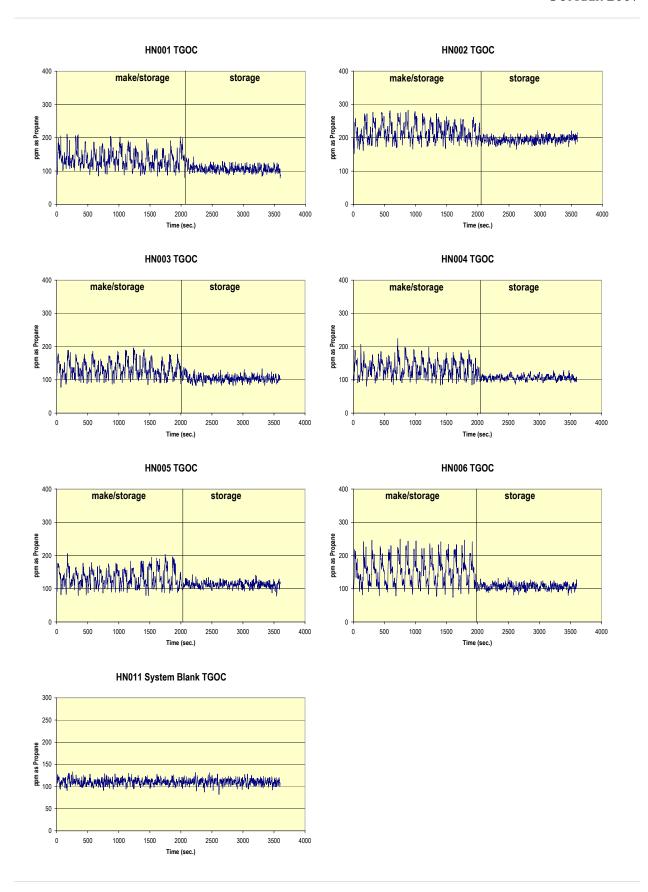
## **Detailed Process Data - Test Series HO**

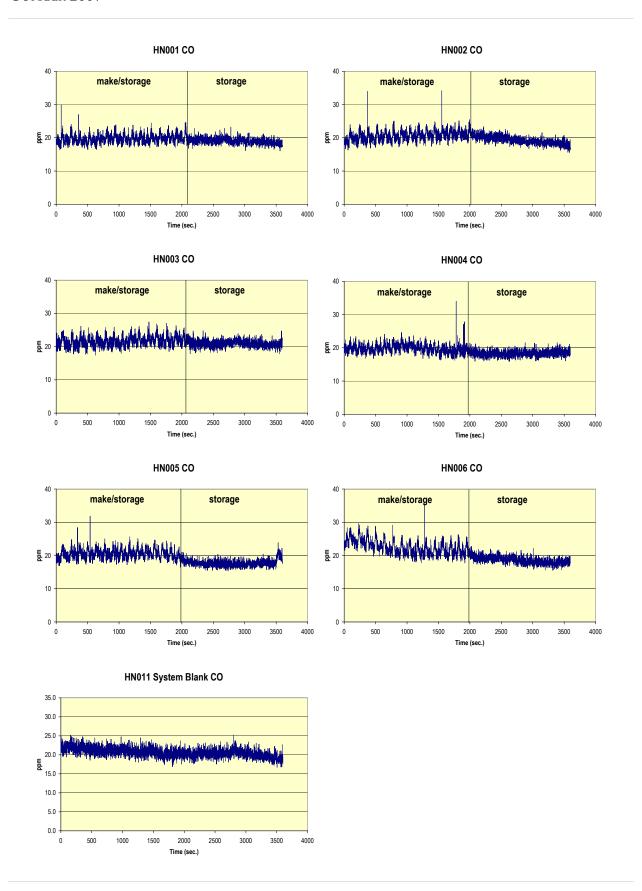
	Technis	and 630BN G	iold Coremak	king			
Test Dates	04/25/07	04/25/07	04/26/07	04/26/07	04/26/07	04/26/07	
Emissions Sample #	HO001	HO002	HO003	HO004	HO005	HO006	Averages
Production Sample #	HO001	HO002	HO003	HO004	HO005	HO006	
Average Cooled Core Weight, g	2570.6	2646.2	2567.1	2655.9	2634.5	2633.2	2617.9
Average Core Box Temperature	439	448	453	455	450	444	448
% stated core binder (BOS)	3	3	3	3	3	3	3
% calculated actual binder	2.9	2.9	2.9	2.9	2.9	2.9	2.9
Number of cores made in run	15	15	15	15	15	15	15
Total core weight in from run, Lbs.	85.01	87.51	84.89	87.83	87.12	87.08	86.57
Total binder weight in run, Lbs.	2.5	2.5	2.5	2.6	2.5	2.5	2.5
Core LOI, %	2.9006	2.8549	ND	2.8598	2.8778	2.8290	2.86

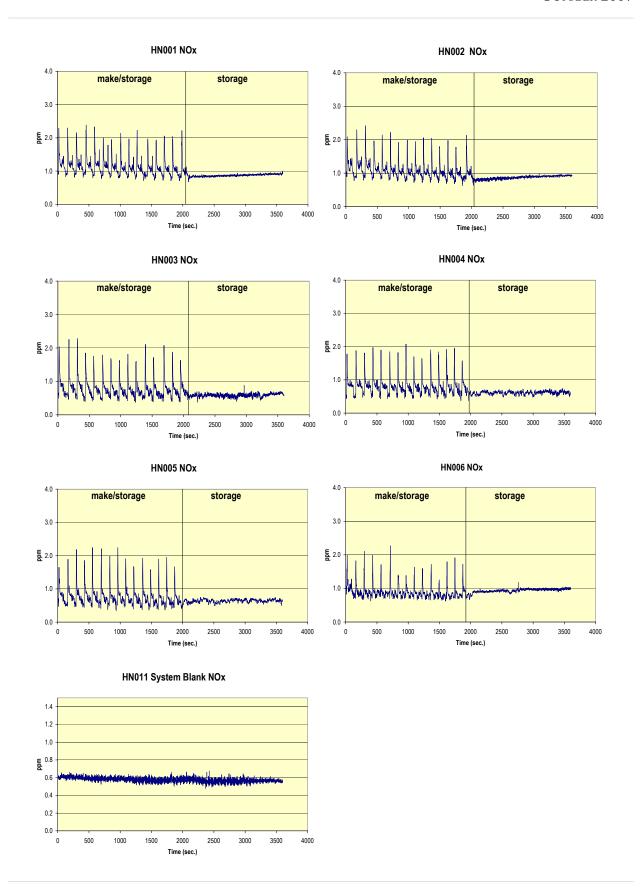


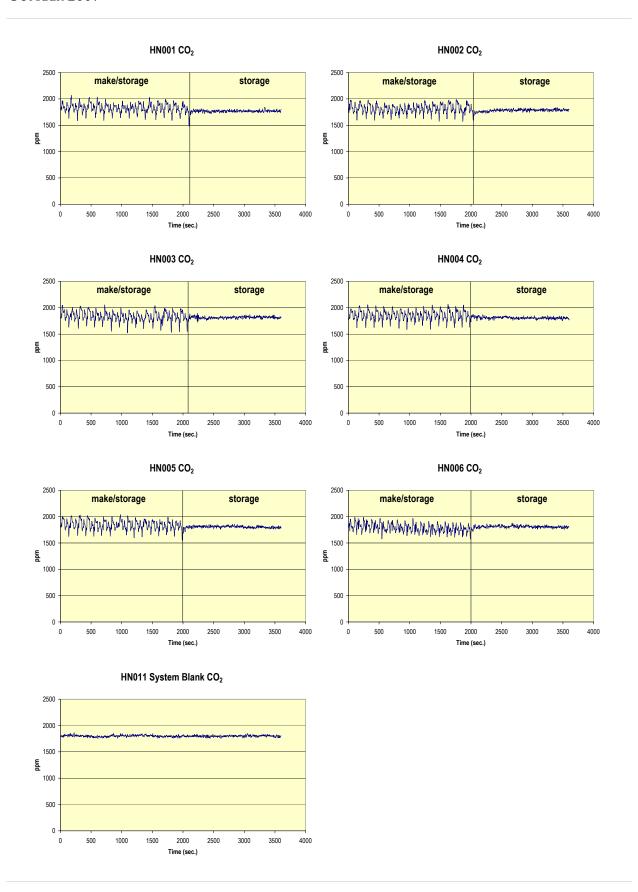
## APPENDIX D CONTINUOUS EMISSION CHARTS

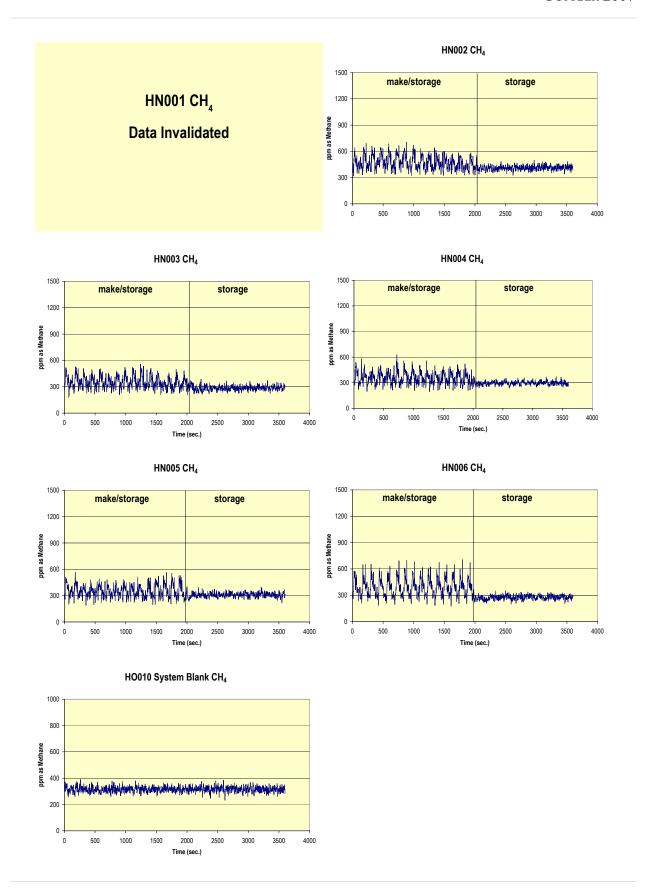


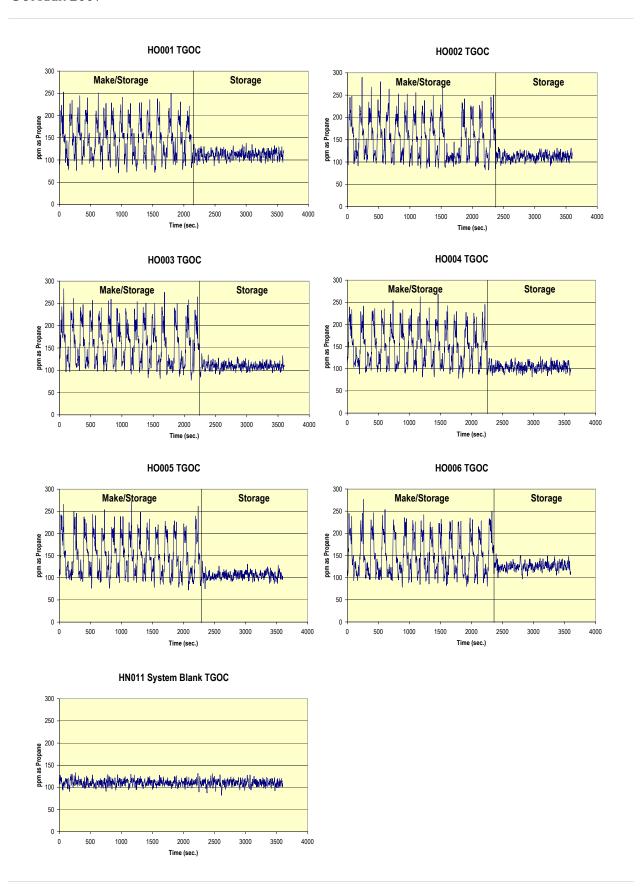


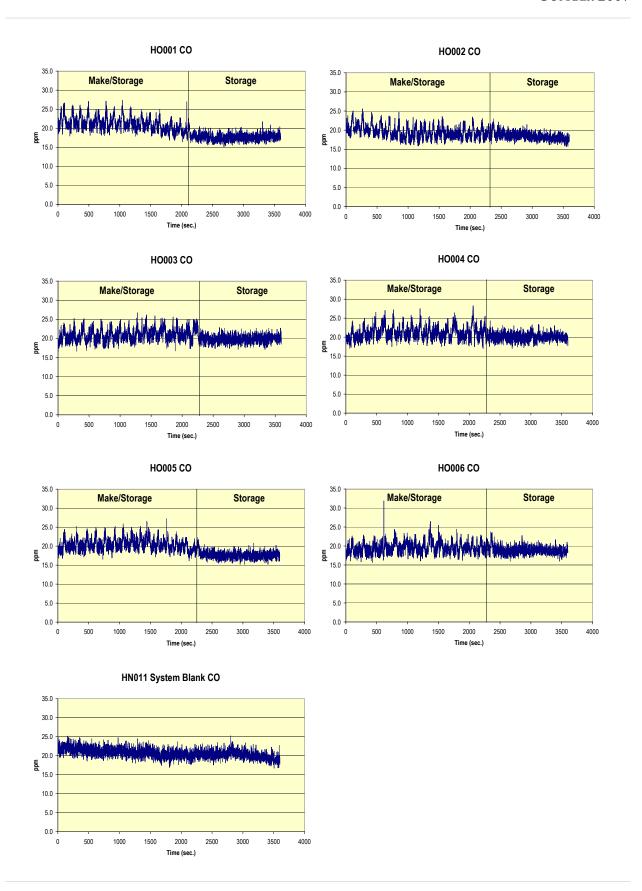


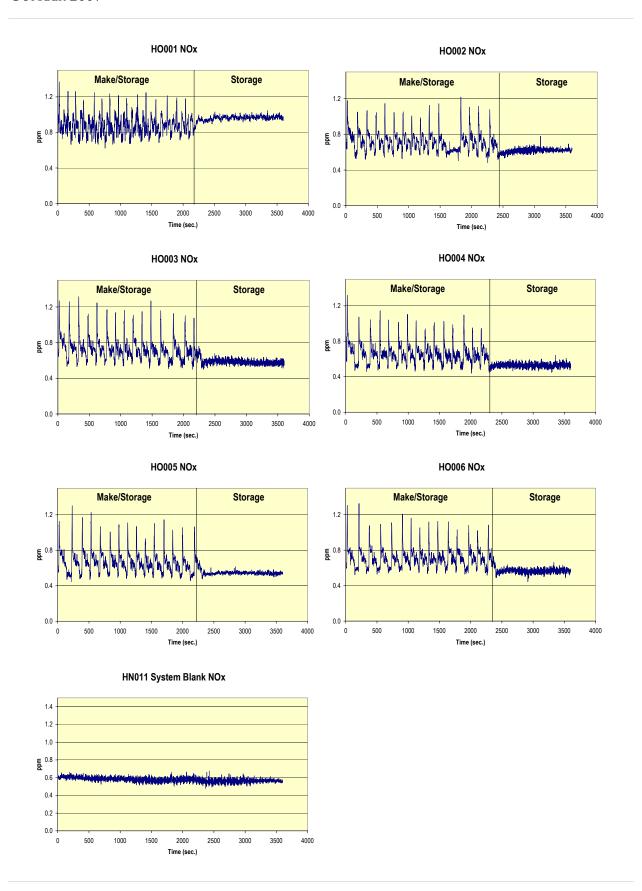


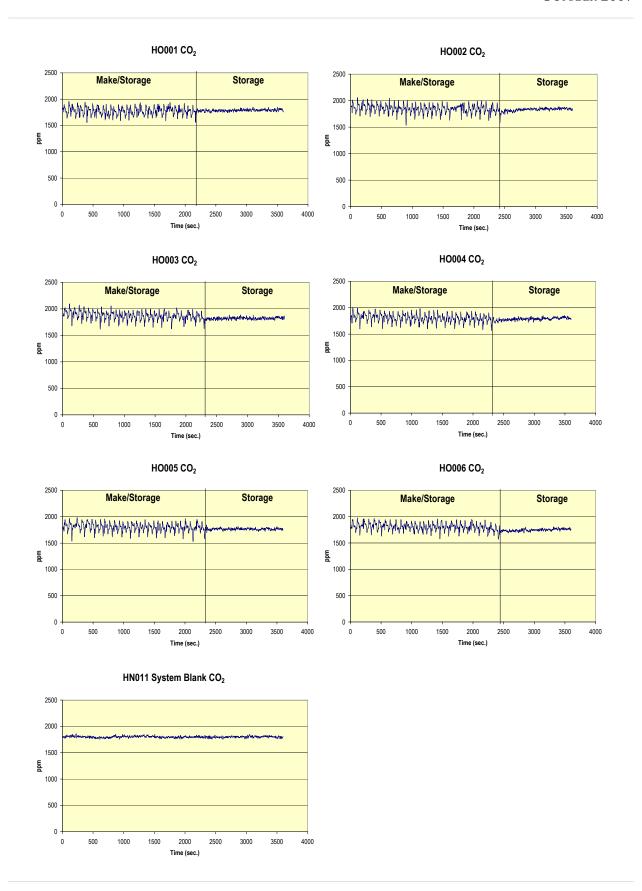


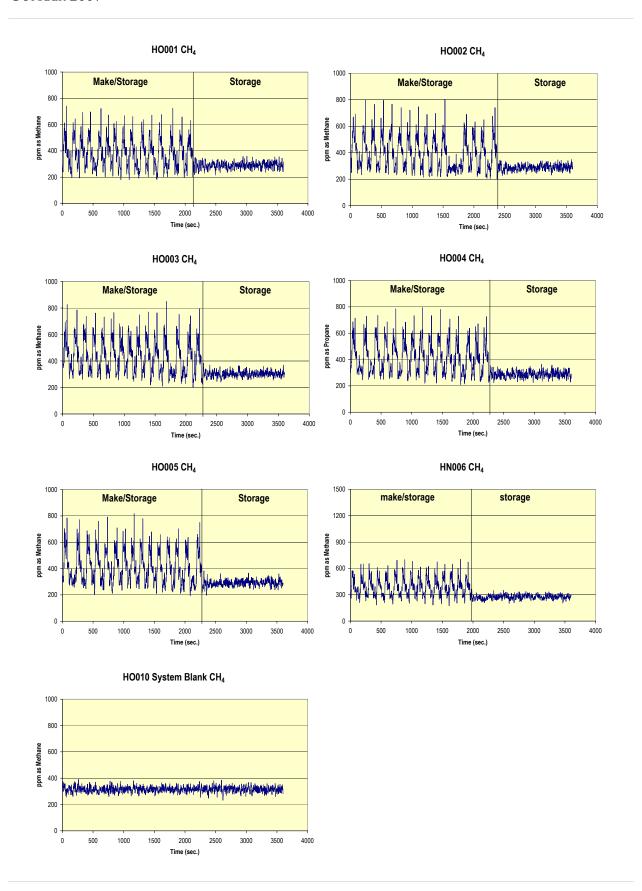


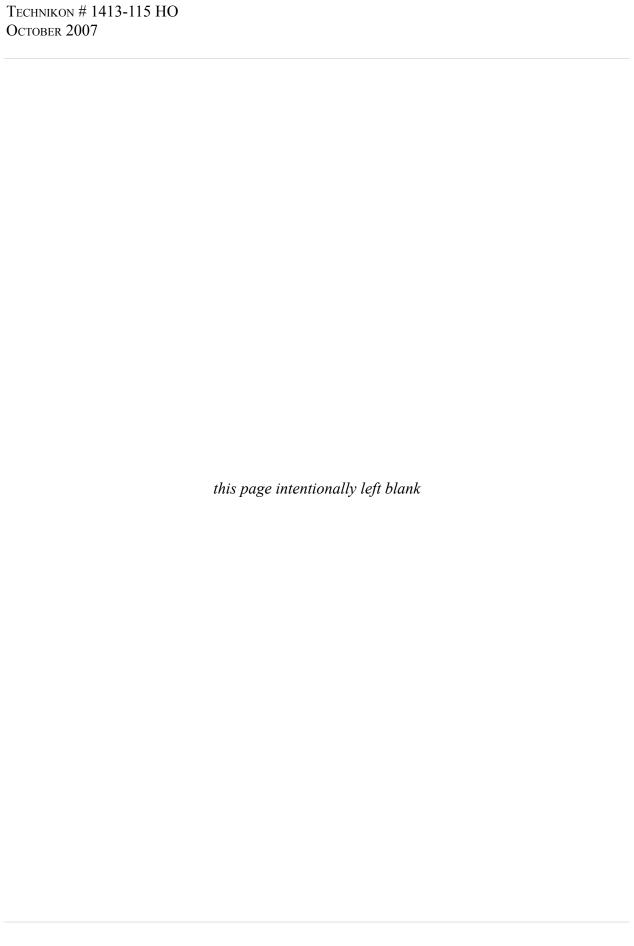












## **Acronyms & Abbreviations**

**AFS** American Foundry Society

**ARDEC** (US) Army Armament Research, Development and Engineering Center

BOS Based on ( ).
BOS Based on Sand.

CAAA Clean Air Act Amendments of 1990CARB California Air Resources Board

**CERP** Casting Emission Reduction Program

**CFR** Code of Federal Regulations

CH<sub>4</sub> Methane

CISA Casting Industry Suppliers Association

CO Carbon Monoxide CO, Carbon Dioxide

**CRADA** Cooperative Research and Development Agreement

CTM Conditional Test MethodDOD Department of DefenseDOE Department of Energy

EPA Environmental Protection Agency
ERC Environmental Research Consortium

**FID** Flame Ionization Detector

GC/MS Gas Chromatography/Mass Spectrometry

**GS** Greensand

HAP Hazardous Air Pollutant defined by the 1990 Clean Air Act Amendment

HC HydrocarbonI Invalidated Data

Lb/Lb Pound per Pound of Binder used Lb/Tn Pound per Ton of Metal poured

LOI Loss on Ignition
MB Methylene Blue

NA Not Applicable; Not Available

ND Non-Detect; Not detected below the practical quantitation limit

NG Natural Gas

NMHC Non-Methane Hydrocarbons

## TECHNIKON # 1413-115 HO

## October 2007

**NOx** Oxides of Nitrogen

NT Not Tested

PCS Pouring, Cooling, Shakeout
POM Polycyclic Organic Matter
psi Pounds per Square Inch

QA/QC Quality Assurance/Quality Control scfm Standard Cubic Feet per Minute

SO<sub>2</sub> Sulfur DioxideTA Target AnalyteTEA Triethylamine

TGOC Total Gaseous Organic Concentration
THC Total Hydrocarbon Concentration

US EPA United States Environmental Protection Agency
USCAR United States Council for Automotive Research

VOST Volatile Organic Sampling Train
WBS Work Breakdown Structure