# June 10, 2024 SUMMA CANISTER REPORT-ADAMS CITY HIGH SCHOOL COMMERCE CITY NORTH DENVER COMMUNITY AIR MONITORING NETWORK COMMERCE CITY, COLORADO

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## **TABLE OF CONTENTS**

SE	CTION PAGE
EXE	ECUTIVE SUMMARY3
1.0	INTRODUCTION4
	1.1 Air Monitoring Site Description4
2.0	METHODS7
	2.1 Air Sampling Methods7
	2.2 Screening Health Risk Assessment Methods
3.0	RESULTS11
	3.1 Summary of Air Sampling Results
	3.2 Screening Health Risk Assessment Results
4.0	Uncertainty Evaluation
5.0	Program Changes
LIS	T OF APPENDICES
Α	SAMPLE CHAIN OF CUSTODIES
LIS	T OF TABLES
1-1	CCND MONITORS AND SUMMA CANISTER SAMPLING LOCATIONS6
1-2	SELECTED ANALYTES MEASURED IN SUMMA CANISTERS8
1-3	CM3 – ADAMS CITY HIGH SCHOOL LOCATION PLANNED AND SENSOR - TRIGGERED EVENT SAMPLE CONCENTRATIONS (PPBV)12
1-4	SUMMA CANISTER SCREENING HEALTH RISK ASSESSMENT: COMPOUND- SPECIFIC HAZARD QUOTIENTS AND HAZARD INDICES FOR CCND CM3 – ADAMS CITY HIGH SCHOOL MONITORING SITE
LIS	T OF FIGURES
1-1	MAP OF TEN CCND MONITOR LOCATIONS5
1-2	CM3 VOC AND WIND DIRECTION   June 10, 2024, 8:36 P.M. – 10:36 P.M
1-3	CM3 WIND ROSE   June 10, 2024, 9:36 P.M. – 10:36 P.M
1-4	COMPOUND-SPECIFIC HAZARD QUOTIENTS FOR VOCS DETECTED IN THE June 10, 2024, SENSOR-TRIGGERED EVENT SAMPLE AT CM3 – ADAMS CITY HIGH SCHOOL LOCATION
1-5	HAZARD INDICIES AT THE CCND CM3 – ADAMS CITY HIGH SCHOOL LOCATION FOR PLANNED AND SENSOR TRIGGERED AIR SAMPLES17

#### **EXECUTIVE SUMMARY**

In response to feedback received by Suncor Energy (U.S.A.) Inc. (Suncor) through community engagement conducted in the fall of 2020, Suncor voluntarily committed to developing a continuous, near real-time air monitoring program to gain insight into air quality for neighborhoods in the vicinity of the Suncor refinery in Commerce City, Colorado. Montrose Environmental Group - Air Quality Services, LLC (Montrose) was contracted by Suncor to deploy, operate and maintain the network in the Commerce City and North Denver (CCND) neighborhoods. Air monitoring was accomplished through three separate technical approaches: (1) continuous, near real-time monitoring for the following analytes: carbon monoxide (CO), sulfur dioxide (SO<sub>2</sub>), hydrogen sulfide (H<sub>2</sub>S), nitrogen oxide or nitric oxide (NO), nitrogen dioxide (NO<sub>2</sub>), particulate matter (PM<sub>2.5</sub>) and total volatile organic compounds (VOCs); (2) periodic collection and laboratory analysis for the presence of specific VOCs from 6-liter evacuated stainless steel ("Summa") canisters; and (3) periodic real-time air monitoring throughout neighborhoods using a mobile monitoring van to detect presence of specific VOCs. An "analyte" is a material that a measuring device is designed to detect and measure. It may be a chemical gas, an airborne particle, or other type of material.

Approach number two consists of collection of air data to measure the presence of specific VOCs. This approach has two parts: collection of planned air samples and collection of unplanned, VOC sensor-triggered air samples. Planned air samples were collected across 13 different locations, ten from within the CCND neighborhoods and three from non-CCND locations (urban and rural background), over a 1-hour time period by a field technician, in Q2 2024. VOC sensor-triggered samples are collected automatically when total VOCs are detected at an airborne concentration of 1 part per million (ppm) or higher for 1 minute or longer. This report analyzes the data from a VOC sensor-triggered air sample collected at Adams City High School (CM3) on June 10, 2024.

Health scientists from CTEH, LLC (CTEH®) (a subsidiary company of Montrose) performed a screening-level human health risk assessment based on the data collected by Montrose. A screening-level assessment uses the most health conservative assumptions about exposure and chemical toxicity. This risk assessment was conducted to determine whether measured concentrations of individual or cumulative (combined) VOCs could potentially cause acute (short-term) adverse health effects. The health risk calculations described in this report were performed per federal and state guidance. The risk assessment of the June 10, 2024, one-hour triggered sample at Adams City High School (CM3) resulted in the following overall findings:

- All measured VOCs (individual and cumulative) were below their respective acute health-based reference levels.
- The cumulative acute health risks calculated from the sensor-triggered event sample were consistent with the planned air sample collected at the same location during the previous quarter.
- The measured concentrations during this triggered sample are not expected to cause an appreciable risk of adverse acute health effects, even for sensitive subpopulations.

### 1.0 INTRODUCTION

In response to feedback received by Suncor Energy (U.S.A.) Inc. (Suncor) through community engagement conducted in the fall of 2020, Suncor voluntarily committed to developing a continuous, near real-time air monitoring program to gain insight into air quality for neighborhoods in the vicinity of the Suncor refinery in Commerce City, Colorado. Montrose Environmental Group - Air Quality Services, LLC (Montrose) was contracted by Suncor to deploy, operate and maintain the network in the Commerce City and North Denver (CCND) neighborhoods. Air monitoring was accomplished through three separate technical approaches: (1) continuous, near real-time monitoring for the following analytes: carbon monoxide (CO), sulfur dioxide (SO<sub>2</sub>), hydrogen sulfide (H<sub>2</sub>S), nitrogen oxide or nitric oxide (NO), nitrogen dioxide (NO<sub>2</sub>), particulate matter (PM<sub>2.5</sub>) and total volatile organic compounds (VOCs); (2) periodic collection and laboratory analysis for the presence of specific VOCs from Summa canisters; and (3) periodic real-time air monitoring throughout neighborhoods using a mobile monitoring van to detect presence of specific VOCs. An "analyte" is a material that a measuring device is designed to detect and measure. It may be a chemical gas, an airborne particle, or other type of material.

The objective of this report is to provide results from a sensor-triggered canister sample collected on June 10, 2024, at the Adams City High School (CM3). The measured concentrations for this single sample were compared to established acute (short-term) health-based reference levels and compared to planned samples collected at the same location.

#### 1.1 Air Monitoring Site Description

Ten monitors and Summa canister sampling locations were positioned throughout the CCND neighborhoods within a three-mile radius of the refinery operations. The monitor locations are shown in Figure 1-1 and described in Table 1-1; they were selected based on the following criteria:

- Historical wind pattern data,
  - Proximity to the refinery and non-refinery sources,
  - Existing infrastructure, as well as site access and safety,
  - Community feedback



FIGURE 1-1
MAP OF TEN CCND MONITOR LOCATIONS

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TABLE 1-1
CCND MONITORS AND SUMMA CANISTER SAMPLING LOCATIONS

Location ID	Secondary ID	GPS Coordinates	Distance from Refinery Center (miles)	Cross Streets
	<u> </u>		(	
CM1	Rose Hill Elementary School	39.80164, -104.90882	2.0	E. 58 <sup>th</sup> Ave. & Oneida St., Commerce City
CM2	Suncor Refinery Business Center	39.79623, -104.95727	0.70	Brighton Blvd. & York St., Commerce City
CM3	Adams City High School	39.82736, -104.90193	2.9	E. 72 <sup>nd</sup> Ave. & Quebec Pkwy, Commerce City
CM4	Adams City Middle School	39.82893, -104.93499	1.9	Birch St. & E. 72 <sup>nd</sup> Ave., Commerce City
CM5	Central Elementary School	39.81365, -104.92191	1.7	Holly St. & E 64 <sup>th</sup> Ave., Commerce City
CM6	Focus Points Family Resource Center	39.78436, -104.95663	1.4	Columbine St. & 48 <sup>th</sup> Ave., Denver
CM7	Kearney Middle School	39.80888, -104.91545	1.7	E. 62 <sup>nd</sup> Ave. & Kearney St., Commerce City
CM8	Monroe	39.81560, -104.94503	0.85	Monroe St. & E. 64 <sup>th</sup> Ave., Denver
СМ9	48 <sup>th</sup> and Race	39.78455, -104.96264	1.7	East 48 <sup>th</sup> Ave. & Race St., Denver
CM10	Alsup Elementary School	39.82026, -104.93663	1.3	East 68 <sup>th</sup> Ave. & Birch St., Commerce City

#### 2.0 METHODS

### 2.1 Air Sampling Methods

A VOC sensor-triggered air sample collection occurred at 9:36 p.m. at the CM3 – Adams City High School location on June 10, 2024.

Air sample collection is triggered by the VOC monitors upon detection of 1 ppm or greater total VOCs for a 1-minute average. A triggered sample is collected over a 1-hour period by an Entech Instruments Silonite™ CS1200E Passive Canister Sampler connected to 6-liter chemically inert stainless steel ("Summa") canister. Prior to deployment, the Summa canister was cleaned and blanked for use according to laboratory Standard Operating Procedures (SOP). Air sampling and analysis was conducted in accordance with the Quality Assurance Project Plan (QAPP) available online at www.ccnd-air.com/documents. The triggered canister sample was shipped to Enthalpy Analytical in Durham, North Carolina. The United States Environmental Protection Agency (USEPA) Compendium Method TO-14A "Determination of Volatile Organic Compounds (VOCs) in Ambient Air using Specially Prepared Canisters with Subsequent Analysis by Gas Chromatography" and TO-15 entitled "Determination of Volatile Organic Compounds (VOCs) in Air Collected in Specially Prepared Canisters and Analyzed by Gas Chromatography/Mass Spectrometry (GC/MS)" was followed for both sampling and analysis methodology. A total of 59 analytes were selected for analysis in this assessment based on the typical suite of analytes monitored for in urban and industrial areas and accounting for laboratory analysis capabilities (Table 1-2).

Planned air samples at ten CCND monitoring locations, which were used in this report to compare to the triggered canister data, were collected during the second quarter of 2024 during a time when near real-time VOC monitors indicated total VOC concentrations to be less than the 1-ppm trigger level. The planned samples were collected and analyzed using the same methods as the triggered sample and full results are available in a separate report.

**TABLE 1-2**SELECTED ANALYTES MEASURED IN SUMMA CANISTERS

Ethylene	Isopentane	3-Methylpentane	3-Methylheptane	2,4- Dimethylpentane
Acetylene	1-Pentene	1-Hexene	Nonane	2,3- Dimethylpentane
Ethane	Pentane	1,3-Butadiene	1,2,3- Trimethylbenzene	
Propylene	Isoprene	Heptane	2-Ethyltoluene	1,3,5- Trimethylbenzene
Propane	Trans-2-Pentene	2-Methylhexane	Decane	2,2,4- Trimethylpentane
Isobutane	Cis-2-Pentene	Cis-2-Pentene Toluene		Tetrachloroethene
1-Butene	2,2- Dimethylbutane	3-Methylhexane	m-Diethylbenzene	1,2,4- Trimethylbenzene
Butane	Cyclopentane	Methylcyclohexane	p-Diethylbenzene	Methylcyclopentane
Trans-2-Butene	Cyclohexane	Hexane	Undecane	2,3,4- Trimethylpentane
Cis-2-Butene	2-Methylpentane	2-Methylheptane	Dodecane	2,3-Dimethylbutane
m-/p-Xylenes	o-Xylene	4-Ethyltoluene	Benzene	Carbon disulfide
n-Octane	Isopropylbenzene	n-Propylbenzene	Naphthalene	

#### 2.2 Screening Health Risk Assessment Methods

CTEH® conducted a screening-level public health risk assessment consistent with federal risk assessment guidelines to determine whether the detected concentrations of individual or cumulative (combined) analytes in the triggered air sample could potentially pose acute (short-term) health impacts and evaluate the data compared to samples collected during planned non-event conditions. A tiered approach to the risk assessment was used. This approach involves one or more iterative steps (or tiers) being performed in which health risks are calculated and evaluated multiple times. In most cases, risk assessors cannot know exactly the level of chemical exposure experienced by individuals or communities.

The first-tier evaluation of the triggered sample made a health-protective assumption that represents an exposure to a person located at that sampling location for an entire hour during the time the sample was collected. Additionally, the first tier assumes that all analytes measured are exerting an effect on the body in a similar manner, which is rarely the case. If the resulting risk values indicate the lack of likely adverse health effects under these worst-case conditions, then the risk assessment is complete. However, if the risk values suggest a potential for adverse health effects, then a second tier of risk calculations are performed, but this time using more detailed assumptions about exposure that are still simple representations of the real world but are more realistic than the first-tier, worst-case assumptions. Each successive tier represents a more complete characterization of exposure variability and/or uncertainty that requires a corresponding increase in calculation complexity and scientific level of effort.

The first tier of this risk assessment process is called a screening-level risk assessment. The conservative assumptions used for this level of risk calculation represent exposure conditions at the distinct sampling location for the entire sampling duration. An exceedance of an acceptable risk level (defined below) does not necessarily indicate that adverse health effects are likely. The Agency for Toxic Substances and Disease Registry (ATSDR) states, "when health assessors find exposures higher than the MRLs (ATSDR's specific health-based reference levels), it means that they may want to look more closely at a site". In other words, screening-level findings of an estimated exposure to a VOC being higher than a health-based reference level do NOT indicate an actual likelihood of adverse effects but do indicate a need to move to a second tier of analysis and refine the risk assessment process with more realistic detail to determine if an actual risk exists that needs to be mitigated.

The screening-level risk assessment reported here includes calculated hazards from exposure to individually measured chemicals as well as exposure to all measured chemicals at once (cumulative). For individual chemicals, an acute health hazard value was calculated as the exposure concentration (EC) divided by the chemical-specific federal or state established human health-based Reference Level (RL) (Equation 1). The result is referred to as the hazard quotient (HQ). Estimates of EC were derived from the 1-hour average concentrations of each analyte. Using the measured values for the EC conservatively assumes that a hypothetical exposed individual occupies the sampling location area and breathes the measured concentration continuously for an hour up to multiple days (an acute exposure).

<sup>&</sup>lt;sup>1</sup>Available at:

https://www.atsdr.cdc.gov/minimalrisklevels/#:~:text=The%20ATSDR%2C%20in%20response%20to,minimal%20risk%20levels%20 (MRLs).

The RLs used to calculate the HQs are previously established exposure levels below which no adverse effect in humans is expected. If available, RLs adopted by the Colorado Department of Public Health and Environment (CDPHE) were selected for use within this assessment. If the analyte was not listed by CDPHE, CTEH® followed a federal and state recommended hierarchy for selection of RLs². Acute HQs were calculated as follows:

#### Eq. 1 – Hazard Quotient (HQ) Equation

HQ = EC / RL

Where:

HQ = Hazard Quotient

EC = 1-hour average air concentration

RL= Acute Health-based Reference Level (from USEPA, ATSDR, Cal EPA and TCEQ)

Health risks from potential cumulative exposures to all detected analytes were calculated by adding together each individual analyte's HQ calculated for a given sampling location. The sum of all the individual HQs is called a Hazard Index (HI). Adding together all the HQs is also a very health-conservative approach because it assumes that all the measured analytes exert an adverse effect on the body in a similar manner, which is rarely the case.

A HQ or HI of less than or equal to one is an indication that the estimated exposure is likely to be without an appreciable risk of adverse health effects, even for sensitive sub-populations. The potential for adverse health effects increases as HQ or HI increase above one, but it is not known by how much. Therefore, calculated hazard values in this assessment that are equal to or less than one indicates an acceptable risk level. HQ or HI values of greater than one would prompt a second-tier risk assessment beyond the screening-level assessment.

According to the USEPA and ATSDR, the federal agencies that establish the RLs note that these values "are set below levels that, based on current information, might cause adverse health effects in the people most sensitive." This is because RLs are based on observed toxicity in human or animal studies with an added safety factor to account for uncertainties in the toxicity data. For example, ATSDR identified the lowest observed adverse effect level (LOAEL) for acute exposure to benzene as 10,200 parts per billion (ppb), based on a study of mice exposed six hours per day for six days. ATSDR then applied a combined safety factor of 300 to derive the final RL to account for several uncertainties, including differences between mice and humans and for sensitive individuals<sup>3</sup>. Therefore, it is scientifically incorrect to assume that all real-world exposures to a chemical at levels higher than an RL likely will result in an adverse effect.

The USEPA also has established values for use in emergency situations, termed Acute Exposure Guideline Levels (AEGLs). Unlike health-based reference levels that can be thousands of times below exposure levels where adverse effects are observed, AEGL values are levels at which different acute adverse health effects may be anticipated to occur. According to USEPA, "AEGL-1 represent exposure levels that could produce mild and progressively increasing but transient and non-disabling odor, taste and sensory irritation or certain asymptomatic, non-sensory effects. With increasing airborne concentration above each AEGL, there is a progressive increase in the

<sup>&</sup>lt;sup>2</sup> Available at: https://drive.google.com/file/d/1P2KEvu0MFiyzQAOQtjQUclqR-WGh1bEX/view

<sup>&</sup>lt;sup>3</sup> Available at: https://www.atsdr.cdc.gov/toxprofiles/tp3-c3.pdf

likelihood of occurrence and the severity of effects described for each corresponding AEGL [i.e., AEGL-2 or AEGL-3]." The AEGL-1 60-minute value, if available for the applicable analyte, was also used for comparison purposes because it is more precautionary (than AEGL-2 or AEGL-3) as the AEGL-1 level reflects potential health impacts that are reversible upon cessation of exposure.

#### 3.0 RESULTS

### 3.1 Summary of Air Sampling Results

The total VOC reading on the CM3 – Adams City High School monitor was part of an event (total VOCs measured above 1 ppm) that occurred at 9:36 pm on June 10, 2024. The total VOC reading was observed above 1 ppm for several minutes and resulted in triggering a 1-hour Summa canister collection. The maximum total VOC concentration reading above 1 ppm during the several minutes was 3.4 ppm. After further investigation onsite, the cause of the Summa trigger could not be determined

The Summa canister's compound-specific concentration results are shown in Table 1-3. Prior to, during and after the total VOC reading above 1 ppm the winds were primarily coming from the South (S) (Figures 1-2 and 1-3). Figure 1-2 provides the 1-minute total VOC concentrations and the wind direction data prior to, during and after this event period. Figure 1-3 displays a wind rose of data collected at the CM3 location from 9:36 p.m. to 10:36 p.m. on June 10, 2024.

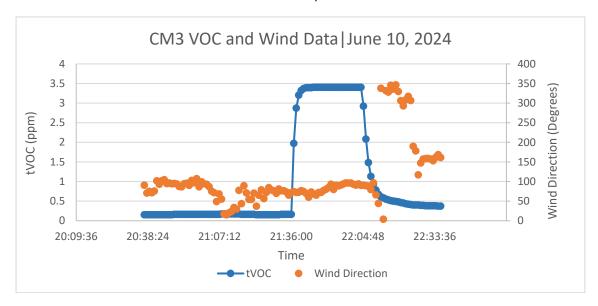
Planned samples at ten CCND sampling locations (including CM3 – Adams high school) were collected in Q2 2024 to evaluate typical VOC levels in the CCND neighborhoods. For comparison, a summary of the planned air sample taken at the CCND CM3 – Adams City High School monitoring location is shown in Table 1-3.

**TABLE 1-3** CM3 - ADAMS CITY HIGH SCHOOL LOCATION PLANNED AND SENSOR -TRIGGERED **EVENT SAMPLE CONCENTRATIONS (PPBV)** 

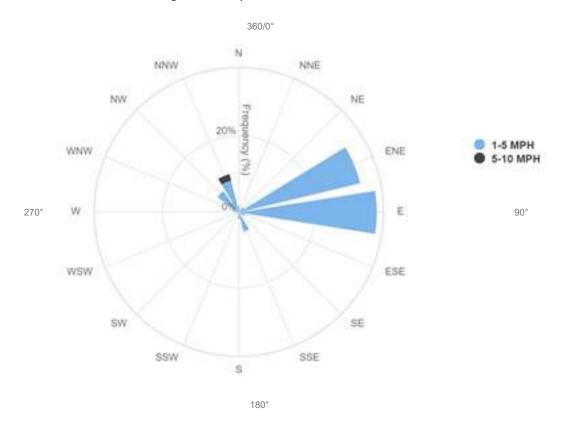
Compound Name			Planned Air Sample	Sensor Triggered Event Sample
1-Butene	Compound Name	Cas No	5/17/2024	6/10/2024
1-Hexnen		106-98-9		
1.   1.   1.   1.   1.   1.   1.   1.				
1.2.3-Trimethylbenzene         \$56.3-6         0.08 (/)         <0.04           1.3-Bitatdiene         106.99-0         <0.04         <0.05           1.3-Diethylbenzene         141.93-5         0.09 (/)         0.11 (/)           1.3-Diethylbenzene         108-67-8         <0.03         <0.04           1.4-Diethylbenzene         105-05-5         0.10 (/)         0.15 (/)           2-Hethylheptane         592-7-8         <0.04         <0.05           2-Methylpetane         592-7-8         <0.04         <0.05           2-Methylpetane         591-76-4         0.10 (/)         0.07 (/)           2-Methylpentane         591-76-4         0.10 (/)         0.07 (/)           2,2-Dimethylbutane         75-83-2         <0.04         <0.05           2,2-Trimethylpentane         56-58-3         <0.04         <0.05           2,3-Dimethylpentane         565-59-3         <0.05 (/)         <0.05           2,3-Dimethylpentane         565-75-3         <0.04         <0.05           2,3-Dimethylpentane         565-75-3         <0.04         <0.05           2,3-Dimethylpentane         565-75-3         <0.04         <0.05           2,4-Dimethylpentane         565-75-3         <0.04         <0.05				
1,2,4-Trimethylbenzene         95-63-6         0.06(j)         < 0.04           1,3-Butadiene         106-99-0         < 0.04         < 0.05           1,3-Diethylbenzene         141-93-5         0.09(j)         0.11(j)           1,3-Erthytolene         105-05-5         0.10(j)         0.15(j,b)           2-Ethytoluene         105-05-5         0.10(j)         0.15(j,b)           2-Methylheptane         592-27-8         < 0.04         < 0.05           2-Methylpentane         107-83-5         0.19(j)         0.07(j)           2-Methylpentane         107-83-5         0.19(j)         0.07(j)           2-Methylpentane         75-83-2         < 0.04         < 0.05           2,2-Dimethylpentane         75-83-2         < 0.04         < 0.05           2,3-Dimethylpentane         75-83-3         0.05(j)         < 0.05           2,3-Dimethylpentane         55-55-3         0.05(j)         < 0.05           2,3-Dimethylpentane         565-55-3         0.05(j)         < 0.05           2,3-Dimethylpentane         565-55-3         0.05(j)         < 0.05           2,3-Dimethylpentane         59-63-6         0.04(j)         < 0.05           2,3-Dimethylpentane         59-65-57-3         0.05(j) <td< th=""><th></th><th></th><th></th><th></th></td<>				
1,3-Bithylberzene   106-99-0   < 0.04   < 0.05   1,3-Dithylberzene   141-93-5   0.09 (1)   0.11 (1)   1,3-Dithylberzene   108-67-8   < 0.03   < 0.04   1,4-Dithylberzene   105-05-5   0.10 (1)   0.15 (1,1b)   0.1				
1,3-5 Tierhylbenzene 141-93-5 0.09 (J) 0.11 (J) 1,3-5 Tierhylbenzene 108-67-8 < 0.03				
1,3,5-rrimethylbenzene   108-67-8   0,00   0,15 (Jb)   0.15 (Jb)				
1.4-Diethylbenzene 105-05-5 0.10 (J) 0.15 (J,b) 2.Hethylheptane 592-27-8				
2-Ethytkoluene 611-14-3 < 0.04				
2-Methylhexane 592-27-8				0.13(3,0)
2-Methylpentane 591-76-4 0.10 (J) 0.07 (J) 2.2-Dimethylputane 107-83-5 0.19 (J) 0.07 (J) 2.2-Dimethylputane 75-83-2 < 0.04 < 0.05 2.2.4-Trimethylpentane 540-84-1 0.10 < 0.04 < 0.05 2.2.4-Trimethylpentane 79-29-8 < 0.04 < 0.05 < 0.05 < 0.04 < 0.05 < 0.05 < 0.04 < 0.05 < 0.05 < 0.04 < 0.05 < 0.05 < 0.04 < 0.05 < 0.05 < 0.05 < 0.04 < 0.05 < 0.05 < 0.04 < 0.05 < 0.05 < 0.05 < 0.04 < 0.05 < 0.05 < 0.05 < 0.04 < 0.05 < 0.05 < 0.05 < 0.05 < 0.04 < 0.05 < 0.05 < 0.05 < 0.04 < 0.05 < 0.05 < 0.05 < 0.04 < 0.05 < 0.05 < 0.05 < 0.04 < 0.05 < 0.05 < 0.05 < 0.04 < 0.05 < 0.05 < 0.05 < 0.04 < 0.05 < 0.05 < 0.04 < 0.05 < 0.05 < 0.04 < 0.05 < 0.05 < 0.04 < 0.05 < 0.05 < 0.04 < 0.05 < 0.05 < 0.04 < 0.05 < 0.05 < 0.04 < 0.05 < 0.05 < 0.04 < 0.05 < 0.05 < 0.04 < 0.05 < 0.05 < 0.04 < 0.05 < 0.05 < 0.04 < 0.05 < 0.05 < 0.04 < 0.05 < 0.05 < 0.04 < 0.05 < 0.05 < 0.04 < 0.05 < 0.05 < 0.04 < 0.05 < 0.05 < 0.04 < 0.05 < 0.05 < 0.04 < 0.05 < 0.05 < 0.04 < 0.05 < 0.05 < 0.04 < 0.05 < 0.05 < 0.04 < 0.05 < 0.05 < 0.04 < 0.05 < 0.05 < 0.04 < 0.05 < 0.05 < 0.04 < 0.05 < 0.05 < 0.04 < 0.05 < 0.05 < 0.04 < 0.05 < 0.05 < 0.04 < 0.05 < 0.05 < 0.04 < 0.05 < 0.05 < 0.04 < 0.05 < 0.05 < 0.04 < 0.05 < 0.05 < 0.04 < 0.05 < 0.05 < 0.04 < 0.05 < 0.05 < 0.04 < 0.05 < 0.05 < 0.05 < 0.04 < 0.05 < 0.05 < 0.05 < 0.04 < 0.05 < 0.05 < 0.05 < 0.04 < 0.05 < 0.05 < 0.05 < 0.05 < 0.05 < 0.05 < 0.05 < 0.05 < 0.05 < 0.05 < 0.05 < 0.05 < 0.05 < 0.05 < 0.05 < 0.05 < 0.05 < 0.05 < 0.05 < 0.05 < 0.05 < 0.05 < 0.05 < 0.05 < 0.05 < 0.05 < 0.05 < 0.05 < 0.05 < 0.05 < 0.05 < 0.05 < 0.05 < 0.05 < 0.05 < 0.05 < 0.05 < 0.05 < 0.05 < 0.05 < 0.05 < 0.05 < 0.05 < 0.05 < 0.05 < 0.05 < 0.05 < 0.05 < 0.05 < 0.05 < 0.05 < 0.05 < 0.05 < 0.05 < 0.05 < 0.05 < 0.05 < 0.05 < 0.05 < 0.05 < 0.05 < 0.05 < 0.05 < 0.05 < 0.05 < 0.05 < 0.05 < 0.05 < 0.05 < 0.05 < 0.05 < 0.05 < 0.05 < 0.05 < 0.05 < 0.05 < 0.05 < 0.05 < 0.05 < 0.05 < 0.05 < 0.05 < 0.05 < 0.05 < 0.05 < 0.05 < 0.05 < 0.05 < 0.05 < 0.05 < 0.05 < 0.05 < 0.05 < 0.05 < 0.05 < 0.05 < 0.05 < 0.05 < 0.05 < 0.05 < 0.05 < 0.05 < 0.05 < 0.05				
2-Methylputane				
2,2-Dimethylbutane         75-83-2         < 0.04         < 0.05           2,2-A-Trimethylputane         79-9-8         < 0.04         < 0.05           2,3-Dimethylputane         565-59-3         0.05 ()         < 0.05           2,3-Dimethylputane         565-59-3         0.06 ()         < 0.05           2,3-Dimethylpentane         565-59-3         < 0.04         < 0.05           2,4-Dimethylpentane         108-08-7         0.10 (J)         < 0.05           2,4-Dimethylpentane         108-08-7         0.10 (J)         < 0.05           3-Ethyltoluene         620-14-4         0.09 (B,J)         0.10 (J)           3-Methylpexane         589-34-4         0.18 (J)         0.20 (J)           3-Methylpexane         96-14-0         0.09 (J)         0.07 (J)           3-Methylpexane         96-14-0         0.09 (J)         0.07 (J)           4-Ethyltoluene         622-96-8         < 0.04         < 0.05           Acetylene         74-86-2         0.38 (J)         0.13 (J)           Benzene         74-86-2         0.38 (J)         0.13 (J)           Berzene         75-15-0         0.04 (J)         0.05 (J)           Cis-2-Butene         590-18-1         < 0.04 (J)         0.05 (J) <th></th> <th></th> <th></th> <th></th>				
2,2,4-Trimethylpentane         540.841         0.10         < 0.04           2,3-Dimethylpentane         79-29-8         < 0.04         < 0.05           2,3-Dimethylpentane         565-59-3         0.05 (J)         < 0.05           2,3-A-Trimethylpentane         108-08-7         0.10 (J)         < 0.05           2,4-Dimethylpentane         108-08-7         0.10 (J)         < 0.05           3-Ethyltoluene         620-14-4         0.09 (B,J)         0.10 (J)           3-Methylheptane         589-31-1         < 0.04         0.32 (J)           3-Methylpentane         589-34-4         0.18 (J)         0.20 (J)           3-Methylpentane         622-96-8         < 0.04         < 0.05 (J)           3-Methylpentane         622-96-8         < 0.04         < 0.05           Acetylene         74-86-2         0.38 (J)         0.13 (J)           4-Ethyltoluene         622-96-8         < 0.04         < 0.05           Acetylene         74-86-2         0.38 (J)         0.13 (J)           Butane         106-97-8         0.70         1.20           Carbon disulfide         75-15-0         0.04 (J)         0.05 (J)           Cis-2-Pentene         627-20-3         < 0.04 (J)         0.05 (J) <th></th> <th></th> <th></th> <th></th>				
2.3-Dimethylbutane 79-9-8				
2.3-Dimethylpentane         565-59-3         0.05 (J)         < 0.05           2.3-A-Trimethylpentane         565-75-3         < 0.04         < 0.05           2.4-Dimethylpentane         108-08-7         0.10 (J)         < 0.05           3-Ethyltoluene         620-14-4         0.09 (B,J)         0.10 (J)           3-Methylheptane         589-31-1         < 0.04         0.32 (J)           3-Methylpentane         589-34-4         0.18 (J)         0.20 (J)           3-Methylpentane         622-96-8         < 0.04         < 0.05 (J)           3-Methylpentane         622-96-8         < 0.04         < 0.05 (J)           4-Ethyltoluene         74-86-2         0.38 (J)         0.13 (J)           Benzene         71-43-2         0.17         0.06 (J)           Butane         106-97-8         0.70         1.20           Carbon disulfide         75-15-0         0.04 (J)         0.05 (J)           Cis-2-Pentene         627-20-3         < 0.04 (J)         0.05 (J)				
2.3.4-Trimethylpentane         565-75-3         < 0.04         < 0.05           2.4-Dimethylpentane         108-08-7         0.10 (J)         < 0.05           3-Ethyltoluene         620-14-4         0.09 (B.J)         0.10 (J)           3-Methylhexane         589-81-1         < 0.04         0.32 (J)           3-Methylpentane         96-14-0         0.09 (J)         0.07 (J)           3-Methylpentane         96-14-0         0.09 (J)         0.07 (J)           4-Ethyltoluene         622-96-8         < 0.04         < 0.05           Acetylene         74-86-2         0.38 (J)         0.13 (J)           Benzene         71-43-2         0.17         0.06 (J)           Butane         106-97-8         0.70         1.20           Carbon disulfide         75-15-0         0.04 (J)         0.05 (J)           Cis-2-Butene         590-18-1         < 0.04         < 0.05           Cis-2-Pentene         627-20-3         < 0.04         < 0.05           Cyclohexane         110-82-7         0.07         < 0.05           Cyclopentane         287-92-3         0.11 (J)         < 0.05           Decane         12-40-3         0.06 (J)         0.09 (J)           Dodecane				
2,4-Dimethylpentane         108-08-7         0.10 (J)         <0.05           3-Ethylstoluene         620-14-4         0.09 (B.)         0.10 (J)           3-Methylheptane         589-81-1         <0.04         0.32 (J)           3-Methylpentane         589-34-4         0.18 (J)         0.20 (J)           3-Methylpentane         622-96-8         <0.04         <0.05           Acetylene         74-86-2         0.38 (J)         0.13 (J)           Benzene         71-43-2         0.17         0.06 (J)           Butane         106-97-8         0.70         1.20           Carbon disulfide         75-15-0         0.04 (J)         0.05 (J)           Cis-2-Butene         590-18-1         <0.04         <0.05           Cis-2-Pentene         627-20-3         <0.04         <0.05           Cyclopentane         110-82-7         0.07         <0.05           Cyclopentane         287-92-3         0.11 (J)         <0.05           Cyclopentane         124-18-5         0.06 (J)         0.09 (J)           Dedecane         124-18-5         0.06 (J)         0.09 (J)           Detacene         124-18-5         0.06 (J)         0.09 (J)           Ethylene         74-84-0 <th></th> <th></th> <th></th> <th></th>				
3-Ethyltoluéne 620-14-4 0.09 (B.J) 0.10 (J) 3-Methylheptane 589-81-1 < 0.04 0.32 (J) 3-Methylhexane 589-81-1 < 0.06 (J) 0.20 (J) 3-Methylhexane 96-14-0 0.09 (J) 0.07 (J) 4-Ethyltoluene 96-14-0 0.09 (J) 0.07 (J) 4-Ethyltoluene 74-86-2 0.38 (J) 0.13 (J) Benzene 74-48-2 0.38 (J) 0.13 (J) Benzene 74-43-2 0.17 0.06 (J) Butane 106-97-8 0.70 1.20 Carbon disulfide 75-15-0 0.04 (J) 0.05 (J) Cis-2-Butene 590-18-1 < 0.04 (-0.05 Cis-2-Pentene 627-20-3 0.04 0.05 Cyclohexane 110-82-7 0.07 0.05 Cyclopentane 287-92-3 0.11 (J) 0.05 (J) Cyclopentane 287-92-3 0.11 (J) 0.09 (J) Dodecane 124-18-5 0.06 (J) 0.09 (J) Dodecane 124-18-5 0.06 (J) 0.09 (J) Ethylbenzene 100-41-4 0.09 0.05 Ethylbenzene 74-88-1 0.78 0.47 (J) Heptane 142-82-5 0.10 (J) 0.06 (J) Heyane 110-54-3 0.24 0.13 Isobutane 75-28-5 0.85 0.41 (J) Isopentane 78-78-4 1.40 0.43 (J) Isopentane 78-78-5 0.08 (J) 0.03 (-0.04 Mp-Xylenes 179601-23-1 0.24 0.08 Methylcyclopentane 98-82-8 0.09 (J) 0.06 (J) 0.06 (J) Naphthalene 91-18-82 0.06 (J) 0.07 (J) Naphthalene 91-20-3 0.05 (J) 0.06 (J) Naphthalene 91-20-3 0.05 (J) 0.06 (J) Naphthalene 91-36-6 0.85 0.47 (J) Propane 74-98-6 1.80 2.20 Propylenzene 115-67-1 0.22 (J) 0.11 (J) Tetrachloroethene 127-18-4 0.03 (-0.04 Tetrachloroethene 128-8-3 0.71 0.22 Trans-2-Butene 624-64-6 (-0.04 0.06 (J) Trans-2-Pentene 624-64-6 (-0.04 0.06 (J) Trans-2-Pentene 624-64-6 (-0.04 0.06 (J)				
3-Methylheptane 589-81-1				
3-Methylpentane 589-34-4 0.18 (J) 0.20 (J) 3-Methylpentane 96-14-0 0.09 (J) 0.07 (J) 4-Ethyltoluene 622-96-8 < 0.04 < 0.05 Acetylene 74-86-2 0.38 (J) 0.13 (J) 8enzene 71-43-2 0.17 0.06 (J) 8utane 106-97-8 0.70 1.20 (Carbon disulfide 75-15-0 0.04 (J) 0.05 (J) (Cis-2-Butene 590-18-1 < 0.04 (J) 0.05 (J) (Cis-2-Pentene 627-20-3 < 0.04 (J) 0.05 (J) (Cis-2-Pentene 627-20-3 < 0.04 < 0.05 (Cis-2-Pentene 627-20-3 < 0.04 < 0.05 (Cis-2-Pentene 627-20-3 < 0.04 (J) 0.05 (J) (Cis-2-Pentene 627-20-3 & 0.01 (J) 0.09 (J) 0.05 (J) (J)				
3-Methylpentane 96-14-0 0.09 (j) 0.07 (j) 4-Ethyltolune 622-96-8 < 0.04 < 0.05   Acetylene 74-86-2 0.38 (j) 0.13 (j) 8-18-18-18-18-18-18-18-18-18-18-18-18-18				
4-Ethyltoluene         622-96-8         < 0.04				
Acetylene         74-86-2         0.38 (J)         0.13 (J)           Benzene         71-43-2         0.17         0.06 (J)           Butane         106-97-8         0.70         1.20           Carbon disulfide         75-15-0         0.04 (J)         0.05 (J)           Cis-2-Butene         690-18-1         < 0.04         < 0.05           Cis-2-Pentene         627-20-3         < 0.04         < 0.05           Cyclopentane         110-82-7         0.07         < 0.05           Cyclopentane         287-92-3         0.11 (J)         < 0.05           Decane         124-18-5         0.06 (J)         0.09 (J)           Dodecane         112-40-3         0.06 (J)         0.09 (J)           Detane         124-18-5         0.06 (J)         0.09 (J)           Decane         112-40-3         0.06 (J)         0.09 (J)           Decane         112-40-3         0.06 (J)         0.09 (J)           Ethane         74-84-0         5.90         4.40           Ethylene         14-84-0         5.90         4.40           Ethylene         14-85-1         0.78         0.47 (J)           Heptane         14-82-5         0.10 (J)         0.06 (J) </th <th></th> <th></th> <th></th> <th></th>				
Benzene         71-43-2         0.17         0.06 (J)           Butane         106-97-8         0.70         1.20           Carbon disulfide         75-15-0         0.04 (J)         0.05 (J)           Cis-2-Butene         590-18-1         < 0.04         < 0.05           Cis-2-Pentene         627-20-3         < 0.04         < 0.05           Cyclohexane         110-82-7         0.07         < 0.05           Cyclopentane         287-92-3         0.11 (J)         < 0.05           Decane         124-18-5         0.06 (J)         0.09 (J)           Dodecane         112-40-3         0.06 (J)         0.09 (J)           Dodecane         112-40-3         0.06 (J)         0.09 (J)           Ethylene         74-84-0         5.90         4.40           Ethylene         74-84-1         0.99         < 0.04           Ethylene         74-85-1         0.78         0.47 (J)           Heptane         142-82-5         0.10 (J)         0.06 (J)           Hexane         110-54-3         0.24         0.13           Isopentane         75-28-5         0.85         0.41 (J)           Isoprentane         78-79-5         0.85 (J)         0.06 (J) <th></th> <th></th> <th></th> <th></th>				
Butane         106-97-8         0.70         1.20           Carbon disulfide         75-15-0         0.04 (J)         0.05 (J)           Cis-2-Butene         590-18-1         < 0.04         < 0.05           Cis-2-Pentene         627-20-3         < 0.04         < 0.05           Cyclohexane         110-82-7         0.07         < 0.05           Cyclopertane         287-92-3         0.11 (J)         < 0.05           Decane         124-18-5         0.06 (J)         0.09 (J)           Dodecane         112-40-3         0.06 (J)         0.09 (J)           Dodecane         112-40-3         0.06 (J)         0.23 (J,b)           Ethane         74-84-0         5.90         4.40           Ethylene         74-84-1         0.78         0.47 (J)           Heptane         142-82-5         0.10 (J)         0.06 (J)           Hexane         110-54-3         0.24         0.13           Isobutane         75-28-5         0.85         0.41 (J)           Isoperne         78-78-4         1.40         0.43 (J)           Isopropylbenzene         78-79-5         0.08 (J)         < 0.05           Isopropylbenzene         179-60-123-1         0.24         < 0				
Carbon disulfide         75-15-0         0.04 (J)         0.05 (J)           Cis-2-Butene         590-18-1         < 0.04         < 0.05           Cis-2-Pentene         627-20-3         < 0.04         < 0.05           Cyclopexane         110-82-7         0.07         < 0.05           Cyclopentane         287-92-3         0.11 (J)         < 0.05           Decane         124-18-5         0.06 (J)         0.09 (J)           Dodecane         112-40-3         0.06 (J)         0.09 (J)           Dodecane         112-40-3         0.06 (J)         0.09 (J)           Ethane         74-84-0         5.90         4.40           Ethylene         74-84-1         0.09         < 0.04           Ethylene         74-85-1         0.78         0.47 (J)           Heptane         142-82-5         0.10 (J)         0.06 (J)           Hexane         110-54-3         0.24         0.13           Isopatiane         75-28-5         0.85         0.41 (J)           Isopertane         78-79-5         0.08 (J)         < 0.05           Isopropylbenzene         78-79-5         0.08 (J)         < 0.05           Isopropylbenzene         98-82-8         < 0.03				
Cis-2-Butene         590-18-1         < 0.04         < 0.05           Cis-2-Pentene         627-20-3         < 0.04         < 0.05           Cyclohexane         110-82-7         0.07         < 0.05           Cyclopentane         287-92-3         0.11 (J)         < 0.05           Decane         124-18-5         0.06 (J)         0.09 (J)           Dodecane         112-40-3         0.06 (J,b)         0.23 (J,b)           Ethane         74-84-0         5.90         4.40           Ethylene         10-41-4         0.09         < 0.04           Ethylene         74-85-1         0.78         0.47 (J)           Heptane         142-82-5         0.10 (J)         0.06 (J)           Hexane         110-54-3         0.24         0.13           Isobutane         75-28-5         0.85         0.41 (J)           Isopentane         78-78-4         1.40         0.43 (J)           Isoprene         78-79-5         0.08 (J)         < 0.05           Isopropylbenzene         98-82-8         < 0.03         < 0.04           Mg-tylcplexane         179601-23-1         0.24         < 0.08           Methylcyclopentane         96-37-7         0.09 (J)         0.06				
Cis-2-Pentene         627-20-3         < 0.04				
Cyclohexane         110-82-7         0.07         < 0.05				
Cyclopentane         287-92-3         0.11 (J)         < 0.05				
Décane         124-18-5         0.06 (J)         0.09 (J)           Dodecane         112-40-3         0.06 (J,b)         0.23 (J,b)           Ethane         74-84-0         5.90         4.40           Ethylene         100-41-4         0.09         < 0.04           Ethylene         74-85-1         0.78         0.47 (J)           Heptane         142-82-5         0.10 (J)         0.06 (J)           Hexane         110-54-3         0.24         0.13           Isobutane         75-28-5         0.85         0.41 (J)           Isoprentane         78-78-4         1.40         0.43 (J)           Isoprene         78-78-4         1.40         0.43 (J)           Isopropylbenzene         98-82-8         < 0.03         < 0.05           Isopropylbenzene         98-82-8         < 0.03         < 0.04           m,p-Xylenes         179601-23-1         0.24         < 0.08           Methylcyclohexane         108-87-2         0.11 (J)         < 0.05           Methylcyclopentane         96-37-7         0.09 (J)         0.06 (J)           n-Octane         111-65-9         0.06 (J)         0.07 (J)           Naphthalene         91-20-3         < 0.05         <				
Dodecane         112-40-3         0.06 (J,b)         0.23 (J,b)           Ethane         74-84-0         5.90         4.40           Ethylbenzene         100-41-4         0.09         < 0.04           Ethylene         74-85-1         0.78         0.47 (J)           Heptane         142-82-5         0.10 (J)         0.06 (J)           Hexane         110-54-3         0.24         0.13           Isobutane         75-28-5         0.85         0.41 (J)           Isopentane         78-78-4         1.40         0.43 (J)           Isoprene         78-79-5         0.08 (J)         < 0.05           Isopropylbenzene         98-82-8         < 0.03         < 0.04           m,p-Xylenes         179601-23-1         0.24         < 0.08           Methylcyclopexane         108-87-2         0.11 (J)         < 0.05           Methylcyclopentane         96-37-7         0.09 (J)         0.06 (J)           n-Octane         111-65-9         0.06 (J)         0.07 (J)           Naphthalene         91-20-3         < 0.05         < 0.06           Nonane         111-84-2         < 0.04         0.06 (J)           0-Xylene         95-47-6         0.09         < 0.				
Ethane         74-84-0         5.90         4.40           Ethylbenzene         100-41-4         0.09         < 0.04           Ethylene         74-85-1         0.78         0.47 (J)           Heptane         142-82-5         0.10 (J)         0.06 (J)           Hexane         110-54-3         0.24         0.13           Isobutane         75-28-5         0.85         0.41 (J)           Isopertane         78-78-4         1.40         0.43 (J)           Isopropylbenzene         98-82-8         < 0.03         < 0.05           Isopropylbenzene         98-82-8         < 0.03         < 0.04           Methylcyclohexane         179601-23-1         0.24         < 0.08           Methylcyclopentane         96-37-7         0.09 (J)         0.06 (J)           Nonane         111-65-9         0.06 (J)         0.07 (J)           Naphthalene         91-20-3         < 0.05         < 0.06           Nonane         111-84-2         < 0.04         0.06 (J)           0-Xylene         95-47-6         0.09         < 0.04           Pentane         109-66-0         0.85         0.47 (J)           Propane         74-98-6         1.80         2.20 <th></th> <th></th> <th></th> <th></th>				
Ethylbenzene         100-41-4         0.09         < 0.04				
Ethylene         74-85-1         0.78         0.47 (J)           Heptane         142-82-5         0.10 (J)         0.06 (J)           Hexane         110-54-3         0.24         0.13           Isobutane         75-28-5         0.85         0.41 (J)           Isopentane         78-78-4         1.40         0.43 (J)           Isoprene         78-79-5         0.08 (J)         < 0.05           Isopropylbenzene         98-82-8         < 0.03         < 0.04           mp-Xylenes         179601-23-1         0.24         < 0.08           Methylcyclopexane         108-87-2         0.11 (J)         < 0.05           Methylcyclopentane         96-37-7         0.09 (J)         0.06 (J)           n-Octane         111-65-9         0.06 (J)         0.07 (J)           Naphthalene         91-20-3         < 0.05         < 0.06           Nonane         111-84-2         < 0.04         0.06 (J)           o-Xylene         95-47-6         0.09         < 0.04           Pentane         109-66-0         0.85         0.47 (J)           Propane         74-98-6         1.80         2.20           Propylbenzene         103-65-1         < 0.03         < 0.04 <th></th> <th></th> <th></th> <th></th>				
Heptane         142-82-5         0.10 (J)         0.06 (J)           Hexane         110-54-3         0.24         0.13           Isobutane         75-28-5         0.85         0.41 (J)           Isoprene         78-78-4         1.40         0.43 (J)           Isoprene         78-79-5         0.08 (J)         <0.05           Isopropylbenzene         98-82-8         <0.03         <0.04           Methylcyclohexane         179601-23-1         0.24         <0.08           Methylcyclopentane         96-37-7         0.09 (J)         0.06 (J)           Non-Octane         111-65-9         0.06 (J)         0.07 (J)           Naphthalene         91-20-3         <0.05         <0.06           Nonane         111-84-2         <0.04         0.06 (J)           o-Xylene         95-47-6         0.09         <0.04           Pentane         109-66-0         0.85         0.47 (J)           Propane         74-98-6         1.80         2.20           Propylbenzene         103-65-1         <0.03         <0.04           Propylene         115-07-1         0.22 (J)         0.11 (J)           Tetrachloroethene         127-18-4         <0.03         <0.04				
Hexane				
Isobutane				
Isopentane				
Isoprene         78-79-5         0.08 (J)         < 0.05				
Sopropylbenzene   98-82-8   < 0.03   < 0.04   m.pXylenes   179601-23-1   0.24   < 0.08   Methylcyclohexane   108-87-2   0.11 (J)   < 0.05   Methylcyclopentane   96-37-7   0.09 (J)   0.06 (J)   n-Octane   111-65-9   0.06 (J)   0.07 (J)   Naphthalene   91-20-3   < 0.05   < 0.06   Nonane   111-84-2   < 0.04   0.06 (J)   0.06 (J)   o-Xylene   95-47-6   0.09   < 0.04   O-Xylene   95-47-6   0.09   < 0.04   O-Xylene   109-66-0   0.85   0.47 (J)   O-Xylene   109-66-0   0.85   0.47 (J)   O-Xylene   103-65-1   0.03   < 0.04   O-Xylene   103-65-1   0.22 (J)   0.11 (J)   O-Xylene   108-88-3   0.71   0.22   O-Xylene   108-88-3   0.71   0.22   O-Xylene   108-88-3   0.71   0.22   O-Xylene   0.05   O-Xylene   0				
m,p-Xylenes         179601-23-1         0.24         <0.08           Methylcyclohexane         108-87-2         0.11 (J)         <0.05           Methylcyclopentane         96-37-7         0.09 (J)         0.06 (J)           n-Octane         111-65-9         0.06 (J)         0.07 (J)           Naphthalene         91-20-3         <0.05         <0.06           Nonane         111-84-2         <0.04         0.06 (J)           o-Xylene         95-47-6         0.09         <0.04           Pentane         109-66-0         0.85         0.47 (J)           Propane         74-98-6         1.80         2.20           Propylbenzene         103-65-1         <0.03         <0.04           Propylene         115-07-1         0.22 (J)         0.11 (J)           Tetrachloroethene         127-18-4         <0.03         <0.04           Tolluene         108-88-3         0.71         0.22           Trans-2-Butene         624-64-6         <0.04         <0.05           Trans-2-Pentene         646-04-8         <0.05         <0.06				
Methylcyclohexane         108-87-2         0.11 (J)         < 0.05				
Methylcyclopentane         96-37-7         0.09 (J)         0.06 (J)           n-Octane         111-65-9         0.06 (J)         0.07 (J)           Naphthalene         91-20-3         < 0.05         < 0.06           Nonane         111-84-2         < 0.04         0.06 (J)           o-Xylene         95-47-6         0.09         < 0.04           Pentane         109-66-0         0.85         0.47 (J)           Propane         74-98-6         1.80         2.20           Propylbenzene         103-65-1         < 0.03         < 0.04           Propylene         115-07-1         0.22 (J)         0.11 (J)           Tetrachloroethene         127-18-4         < 0.03         < 0.04           Tolluene         108-88-3         0.71         0.22           Trans-2-Butene         624-64-6         < 0.04         < 0.05           Trans-2-Pentene         646-04-8         < 0.05         < 0.06				
n-Octane         111-65-9         0.06 (J)         0.07 (J)           Naphthalene         91-20-3         < 0.05         < 0.06           Nonane         111-84-2         < 0.04         0.06 (J)           o-Xylene         95-47-6         0.09         < 0.04           Pentane         109-66-0         0.85         0.47 (J)           Propane         74-98-6         1.80         2.20           Propylbenzene         103-65-1         < 0.03         < 0.04           Propylene         115-07-1         0.22 (J)         0.11 (J)           Tetrachloroethene         127-18-4         < 0.03         < 0.04           Tolluene         108-88-3         0.71         0.22           Trans-2-Butene         624-64-6         < 0.04         < 0.05           Trans-2-Pentene         646-04-8         < 0.05         < 0.06				
Naphthalene         91-20-3         < 0.05				
Nonane         111-84-2         < 0.04				
o-Xylene         95-47-6         0.09         < 0.04				
Pentane         109-66-0         0.85         0.47 (J)           Propane         74-98-6         1.80         2.20           Propylbenzene         103-65-1         <0.03         <0.04           Propylene         115-07-1         0.22 (J)         0.11 (J)           Tetrachloroethene         127-18-4         <0.03         <0.04           Toluene         108-88-3         0.71         0.22           Trans-2-Butene         624-64-6         <0.04         <0.05           Trans-2-Pentene         646-04-8         <0.05         <0.06				
Propane         74-98-6         1.80         2.20           Propylbenzene         103-65-1         < 0.03         < 0.04           Propylene         115-07-1         0.22 (J)         0.11 (J)           Tetrachloroethene         127-18-4         < 0.03         < 0.04           Toluene         108-88-3         0.71         0.22           Trans-2-Butene         624-64-6         < 0.04         < 0.05           Trans-2-Pentene         646-04-8         < 0.05         < 0.06				
Propylbenzene         103-65-1         < 0.03				
Propylene         115-07-1         0.22 (J)         0.11 (J)           Tetrachloroethene         127-18-4         < 0.03         < 0.04           Tolluene         108-88-3         0.71         0.22           Trans-2-Butene         624-64-6         < 0.04         < 0.05           Trans-2-Pentene         646-04-8         < 0.05         < 0.06				
Tetrachloroethene         127-18-4         < 0.03				
Toluene         108-88-3         0.71         0.22           Trans-2-Butene         624-64-6         < 0.04	Propylene			
Trans-2-Butene         624-64-6         < 0.04				
<b>Trans-2-Pentene</b> 646-04-8 < 0.05 < 0.06				
		624-64-6		
Undecane 1120-21-4 < 0.04 0.12 (J,b)		646-04-8		
	Undecane	1120-21-4	< 0.04	0.12 (J,b)

All results presented in parts per billion (ppb)
Laboratory non-detections are reported as less than ("<") the method detection limit.
Result qualifiers: (J) flag indicates the reported value is an estimate and was detected below the reporting limit; (B) flag indicates that contamination was found in associated Method Blank; (b) High responses were observed for n-undecane, n-dodecane, and 1,4 diethylbenzene in the ICV analyzed 6/18/24 03:26; these analytes were not detected at or above the RL in the associated samples

**FIGURE 1-2**CM3 TOTAL VOC AND WIND DIRECTION | June 10, 2024, 8:36 P.M. – 10:36 P.M.



**FIGURE 1-3** CM3 Adams High School | June 10 , 2024, 9:36 P.M. – 10:36 P.M.



#### 3.2 Screening Health Risk Assessment Results

The purpose of this screening health risk assessment was to determine whether exposure to the concentrations of individual or cumulative VOCs measured in the June 10, 2024, sensor-triggered event sample, collected at Adams City High School, could potentially pose acute (short-term) health hazards. According to USEPA guidelines (USEPA 1989, 2004), a HQ or HI less than or equal to one indicates that exposures are likely to be without any appreciable risk of adverse acute health effects, even for sensitive sub-populations. The calculated acute HQ and HI are summarized in Table 1-4. In general, the data and health risk assessment indicate:

- The June 10, 2024 sensor-triggered event sample concentrations were below their respective acute health-based reference levels (Table 1-4, Figure 1-4)
- The June 10, 2024, sensor-triggered event sample cumulative hazard index (CM3 HI = 0.02) was consistent with the planned air sample collected at the same location during the same quarter(HI = 0.03, Figure 1-5).
- The measured concentrations during this triggered sample are not expected to cause an appreciable risk of adverse acute health effects, even for sensitive subpopulations.

TABLE 1-4
SUMMA CANISTER SCREENING HEALTH RISK ASSESSMENT: COMPOUND-SPECIFIC HAZARD QUOTIENTS AND HAZARD INDICES FOR CCND CM3 – ADAMS CITY HIGH SCHOOL MONITORING SITE

					Planned Air Sample	Sensor Triggered Event Sample
		AEGL 1 60 min Value	Health Based Reference		May 17, 2024	June 10, 2024
Compound Name	Cas No	(ppb)	Level (ppb)	Source		
1-Butene	106-98-9	NR	27,000	TCEQ Short-Term AMCV	0.0000	0.0000
1-Hexene	592-41-6	NR	500	TCEQ Short-Term AMCV	0.0001	0.0001
1-Pentene	109-67-1	NR	12,000	TCEQ Short-Term AMCV	0.0000	0.0000
	526-73-8	140,000	3,000	TCEQ Short-Term AMCV	0.0000	0.0000
	95-63-6	140,000	3,000 298	TCEQ Short-Term AMCV	0.0000 0.0001	0.0000
1,3-Butadiene 1,3-Diethylbenzene	106-99-0 141-93-5	670,000 NR	450	OEHHA Acute REL TCEO Short-Term AMCV	0.0001	0.0002
1,3,5-Trimethylbenzene		140.000	3,000	TCEQ Short-Term AMCV	0.0002	0.0002
1,4-Diethylbenzene	105-05-5	NR	450	TCEQ Short-Term AMCV	0.0002	0.0003
2-Ethyltoluene	611-14-3	NR	250	TCEQ Short-Term AMCV	0.0002	0.0004
2-Methylheptane	592-27-8	NR	4,100	TCEQ Short-Term AMCV	0.0000	0.0000
2-Methylhexane	591-76-4	NR	8,300	TCEQ Short-Term AMCV	0.0000	0.0000
2-Methylpentane	107-83-5	NR	5,400	TCEQ Short-Term AMCV	0.0000	0.0000
2,2-Dimethylbutane	75-83-2	NR	5,400	TCEQ Short-Term AMCV	0.0000	0.0000
	540-84-1	NR	4,100	TCEQ Short-Term AMCV	0.0000	0.0000
2,3-Dimethylbutane	79-29-8	NR	5,400	TCEQ Short-Term AMCV	0.0000	0.0000
2,3-Dimethylpentane	565-59-3	NR	8,300	TCEQ Short-Term AMCV	0.0000	0.0000
2,3,4-Trimethylpentane 2,4-Dimethylpentane	565-75-3 108-08-7	NR NR	4,100 8,300	TCEQ Short-Term AMCV TCEQ Short-Term AMCV	0.0000	0.0000
3-Ethyltoluene	620-14-4	NR	250	TCEQ Short-Term AMCV	0.0004	0.0004
3-Methylheptane	589-81-1	NR	4,100	TCEQ Short-Term AMCV	0.0004	0.0004
3-Methylhexane	589-34-4	NR	8,300	TCEQ Short-Term AMCV	0.0000	0.0000
3-Methylpentane	96-14-0	NR	5,400	TCEQ Short-Term AMCV	0.0000	0.0000
4-Ethyltoluene	622-96-8	NR	250	TCEQ Short-Term AMCV	0.0001	0.0002
Acetylene	74-86-2	NR	25,000	TCEQ Short-Term AMCV	0.0000	0.0000
Benzene	71-43-2	52,000	9	ATSDR Acute MRL	0.0189	0.0069
Butane	106-97-8	5,500,000		TCEQ Short-Term AMCV	0.0000	0.0000
Carbon disulfide	75-15-0	13,000	1,990	OEHHA Acute REL	0.0000	0.0000
Cis-2-Butene	590-18-1	NR	15,000	TCEQ Short-Term AMCV	0.0000	0.0000
Cis-2-Pentene	627-20-3	NR	12,000	TCEQ Short-Term AMCV	0.0000	0.0000
Cyclohexane Cyclopentane	110-82-7 287-92-3	NR NR	1,000 5,900	TCEQ Short-Term AMCV TCEQ Short-Term AMCV	0.0001	0.0001
Decane	124-18-5	NR	1.000	TCEQ Short-Term AMCV	0.0000	0.0000
Dodecane	112-40-3	NR	1,720	CDPHE Acute	0.0000	0.0001
Ethane	74-84-0	NR	NA NA	NA	NA NA	NA
Ethylbenzene	100-41-4	33,000	5,000	ATSDR Acute MRL	0.0000	0.0000
Ethylene	74-85-1	NŔ	500,000	TCEQ Short-Term AMCV	0.0000	0.0000
Heptane	142-82-5	NR	8,300	TCEQ Short-Term AMCV	0.0000	0.0000
Hexane	110-54-3	NR	5,400	TCEQ Short-Term AMCV	0.0000	0.0000
Isobutane	75-28-5	NR	33,000	TCEQ Short-Term AMCV	0.0000	0.0000
Isopentane	78-78-4	NR	68,000	TCEQ Short-Term AMCV	0.0000	0.0000
Isoprene	78-79-5	NR FO 000	1,400	TCEQ Short-Term AMCV	0.0001	0.0000
Isopropylbenzene m,p-Xylenes	98-82-8 179601-23	50,000 130,000	510 2,000	TCEQ Short-Term AMCV ATSDR Acute MRL	0.0001 0.0001	0.0001 0.0000
Methylcyclohexane	108-87-2	NR	4,000	TCEQ Short-Term AMCV	0.0001	0.0000
Methylcyclopentane	96-37-7	NR	750	TCEQ Short-Term AMCV	0.0001	0.0001
n-Octane	111-65-9	NR	4,100	TCEQ Short-Term AMCV	0.0000	0.0000
Naphthalene	91-20-3	NR	95	TCEQ Short-Term AMCV	0.0005	0.0006
Nonane	111-84-2	NR	3,000	TCEQ Short-Term AMCV	0.0000	0.0000
o-Xylene	95-47-6	130,000	1,700	ATSDR Acute MRL	0.0001	0.0000
Pentane	109-66-0	NR	68,000	TCEQ Short-Term AMCV	0.0000	0.0000
Propane	74-98-6	5,500,000	NA	NA	NA	NA
Propylbenzene	103-65-1	NR	510	TCEQ Short-Term AMCV	0.0001	0.0001
Propylene	115-07-1	NR	NA	NA	NA	NA
Tetrachloroethene	127-18-4	35,000	6	ATSDR Acute MRL	0.0053	0.0065
Toluene Trans 2 Butons	108-88-3	67,000	2,000	ATSDR Acute MRL	0.0004	0.0001
Trans-2-Butene Trans-2-Pentene	624-64-6 646-04-8	NR NR	15,000 12,000	TCEQ Short-Term AMCV TCEQ Short-Term AMCV	0.0000	0.0000
Undecane	1120-21-4	NR NR	550	TCEQ Short-Term AMCV	0.0001	0.0000
	11CU-C1-4	INK	JJU	TOLIQ SHOLE TELLIL AIVICY	0.0001	0.0002

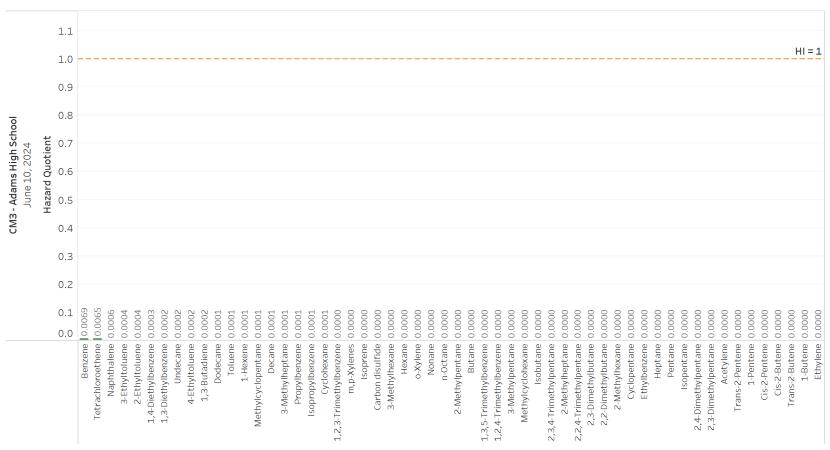
NA = Not Available

NR = According to EPA, AEGL is "Not Recommended due to insufficient data"

pbb - parts per billion

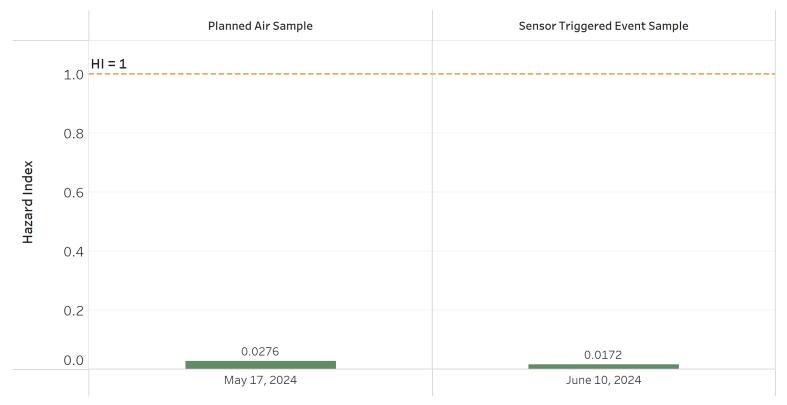
AMCV - Air Monitoring Comparison Value; MRL - Minimum Risk Level; REL - Reference Exposure Level; TCEQ - Texas Commission on Environmental Quality; ATSDR - Agency for Toxic Substances and Disease Registry; OEHHA - Office of Environmental Health Hazard Assessment; CDPHE - Colorado Department of Public Health and Environment

FIGURE 1-4
COMPOUND-SPECIFIC HAZARD QUOTIENTS FOR VOCS DETECTED IN THE June 10, 2024, SENSOR-TRIGGERED EVENT SAMPLE AT CM3 – ADAMS CITY HIGH SCHOOL LOCATION



Hazard quotient (HQ) is the exposure concentration (EC), or air concentration divided by the established health based reference level (RL) for each compound. According to the EPA, a HQ less than 1 (orange line) indicates that exposures are likely to be without appreciable risk of adverse acute health effects, even for sensitive sub-populations. Propylene, propane, and ethane did not have a RL and are not displayed.

FIGURE 1-5
HAZARD INDICIES AT THE CCND CM3 – ADAMS CITY HIGH SCHOOL LOCATION FOR PLANNED AND SENSOR
TRIGGERED AIR SAMPLES



Hazard Index (HI) is the sum of all combined hazard quotients (HQ). According to EPA, a HI less than or equal to one (orange line) indicates that exposures are likely to be without any appreciable risk of adverse acute health effects, even for sensitive sub-populations.

## 4.0 Uncertainty Evaluation

Scientific uncertainty is inherent in each step of the risk assessment process because all risk assessments incorporate a variety of assumptions and professional judgments (USEPA 1989, 2004). Therefore, the acute hazard estimates presented in this assessment are conditional estimates given a considerable number of assumptions about exposure and toxicity. This screening-level risk assessment relied on a combination of health-protective scenarios and toxicity information (i.e., exposure durations, toxicological uncertainty factors, cumulative risk evaluations). This approach was selected to help risk management decision making. Because of these assumptions, the estimates of acute hazards are themselves uncertain but likely to be overestimated compared to actual.

This risk assessment did not address past or present health outcomes associated with current or past exposures. As such, this risk assessment cannot be used to make realistic predictions of biological effects and/or used to determine whether someone is ill (cancer or other adverse health effects) due to past or current exposures. This risk assessment was limited to inhalation exposures from outdoor exposures to all potential sources.

## 5.0 Program Changes

None at this time.

Prepared by:

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Montrose Air Quality Services

Michael Lumpkin, PhD, DABT Senior Toxicologist

Michael H. Lungshin

CTEH®, LLC

## **APPENDIX ASAMPLE CHAIN OF CUSTODIES**

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					Report To:	David	Smith			Number:		PROJ-022555							
Special Instructions:					Email:	davids	mith1@	montrose	env.com	P.O. #:	P.O. #:				PO-065450				
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