

Bristol-Myers Squibb

Building 3 Vapor Intrusion and Air Treatment System Assessment Report

Syracuse North Campus Restoration Area East Syracuse, New York

October 2021

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Acronyms and Abbreviations

Arcadis Arcadis of New York, Inc.

ASHRAE American Society of Heating, Refrigerating and Air-Conditioning Engineers

ATS air treatment system

BASE Building Assessment Survey and Evaluation Study

BCA Brownfield Cleanup Agreement
BDA Brownfield Development Area

BMS Bristol-Myers Squibb

CFM cubic feet per minute
COC constituent of concern

DER-10 DER-10 Technical Guidance for Site Investigation and Remediation

DUSR Data Usability Summary Report

ELAP Environmental Laboratory Approval Program

eV electron volt

FSAP Field Sampling and Analysis Plan

HVAC heating, ventilation, and air conditioning

JLL Jones Lang LaSalle Incorporated

NYSDEC New York State Department of Environmental Conservation

NYSDOH New York State Department of Health

OBG O'Brien & Gere

OMM operation, maintenance, and monitoring

OSHA Occupational Safety and Health Administration

Phase 1/1A Report Phase 1/1A Remedial Investigation Data Summary Report

PID photoionization detector

ppm parts per million

QAPP Quality Assurance Project Plan
QA/QC quality assurance/quality control

RSL Regional Screening Level

SDS Safety Data Sheet
SVI soil vapor intrusion

TIC tentatively identified compound

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TCE trichloroethene

TLV Threshold Limit Values

μg/L micrograms per liter

μg/m³ micrograms per cubic meter

USEPA United States Environmental Protection Agency

VI vapor intrusion

VOC volatile organic compounds

1 Introduction

This Building 3 Vapor Intrusion and Air Treatment System Assessment Report has been developed by Arcadis of New York, Inc. (Arcadis) for the Bristol-Myers Squibb (BMS) Syracuse North Campus Restoration Area, which is also referred to as the Brownfield Development Area (BDA). The BDA is part of the BMS Facility located at 3551 Burnet Avenue in East Syracuse, New York. The location of the BDA and BMS Facility are shown on **Figure 1**. The BDA is identified by the New York State Department of Environmental Conservation (NYSDEC) as Site No. C734138 and is subject to a Brownfield Cleanup Agreement (BCA; October 2011) between BMS and the New York State Department of Environmental Conservation (NYSDEC).

An indoor air treatment system (ATS) was installed in April 2021 as a supplemental mitigation measure for the boiler control room located in the basement of Building 3 at the BMS Facility. The proposed installation of the ATS was described in a March 19, 2021 letter submitted to NYSDEC by Arcadis on behalf of BMS. This report presents the findings of pre-startup and post-startup (i.e., confirmation sampling) indoor air sampling events conducted in Building 3 in January and June 2021, respectively, as well as supplemental air sampling conducted to assess ATS performance. Additionally, a summary of previous soil vapor intrusion (SVI) assessment activities conducted prior to 2021 at the BMS Facility is included in this report.

2 Previous Building 3 Soil Vapor Intrusion Assessments

Based on the presence of dissolved-phase volatile organic compound (VOC) concentrations in groundwater in the BDA, an evaluation of the potential SVI migration pathway was conducted for Building 3 as it was a downgradient or proximal building that remained occupied at the time of the Phase 1A investigation in 2016.

One indoor air sample (with associated outdoor ambient air sample) was collected in the Building 3 basement (i.e., boiler control room) for each sampling event. An elevated water table in proximity to the Building 3 slab elevation precluded collection of sub-slab samples or near building soil vapor samples. Therefore, BMS proposed to use the analytical results from monitoring wells upgradient of Building 3 (BDA-1WT and BDA-1F, shown on **Figure 2**) and compare compounds detected in the groundwater sample results to the compounds identified in the indoor air sample.

The initial SVI investigation consisted of two (2) rounds of sampling. The first round was conducted in January and March 2016. The results of that sampling were documented in the *Phase 1/1A Remedial Investigation Data Summary Report* (Arcadis 2016) (Phase 1/1A Report). In October 2016, BMS informed the NYSDEC and the New York State Department of Health (NYSDOH) that it would conduct a second round of SVI sampling during the following heating season. The second round of SVI sampling was conducted in December 2016 with additional groundwater samples from upgradient monitoring wells BDA-1F and BDA-1WT collected in July and December of 2017.

A summary of the findings from the 2016 sampling events was submitted to the NYSDEC and the NYSDOH on March 8, 2017 in a report titled *Soil Vapor Intrusion (SVI) Module Data Summary Report* (SVI Report, Arcadis 2017). The SVI Report concluded that based on the 2016 sampling results no additional sampling would be needed. In May 2017, the NYSDOH updated the screening values listed in the decision matrices from the *2006 NYSDOH Guidance for Evaluating Soil Vapor Intrusion in the State of New York*. The results of this change were discussed in an October 4, 2017 e-mail to BMS from the NYSDOH, and during an October 10, 2017 meeting between BMS, the NYSDEC, and the NYSDOH. The agencies expressed concern that the Building 3 measured indoor air concentrations of trichloroethene (TCE) at 0.93 and 1.6 micrograms per cubic meter [µg/m³] were greater than the NYSDOH revised indoor air guidance values for monitoring (0.2 µg/m³) and mitigation (1.0 µg/m³) when a sub-slab or soil gas source is detected. During the October 10, 2017 meeting, BMS indicated that additional sampling of indoor air in Building 3 would be conducted following completion of on-going building renovations in Building 3 and the adjoining Building 2, and the onset of the heating season.

The third round of Building 3 indoor and ambient air sampling (one sample each) was conducted on January 17, 2018 after completing the renovation activities but prior to startup of a new heating, ventilation, and air conditioning (HVAC) system. The results identified a slight decrease in the TCE concentration versus the previous rounds with a concentration of $0.77~\mu g/m^3$ and $0.82~\mu g/m^3$ from the parent and duplicate indoor air samples, respectively. However, the results remained above the screening criteria for monitoring of $0.2~\mu g/m^3$ for TCE. As such, in the March 2018 monthly progress report to the NYSDEC and the NYSDOH, BMS provided an additional work scope and schedule, which was approved on May 2, 2018 following a modification request by the NYSDEC/NYSDOH.

The fourth round of Building 3 indoor and ambient air sampling (one sample each) was conducted on January 24, 2019 following the installation of the passive mitigation measures identified in the March 2018 work scope and

May 2, 2018 approval letter, and with the new HVAC system in full operation. Arcadis conducted a building airflow and pressure assessment on January 11, 2019. The evaluation results were detailed in the *Building 3 Vapor Intrusion (VI) Assessment Report* (Arcadis 2019), which indicated a positive pressure in the boiler control room indoor air relative to other locations measured (e.g., drain feature, indoor air in adjoining building rooms and outdoor ambient air), suggesting conditions suitable for preventing vapors from migrating through the floor drains or from the sub-slab environment and into the breathing zone. The results again identified a slight decrease in the TCE concentration versus the previous rounds with a concentration of 0.59 μ g/m³ and 0.61 μ g/m³ from the parent and duplicate indoor air samples, respectively. However, the results remained above the screening criteria for monitoring of 0.2 μ g/m³ for TCE. As such, BMS proposed to the NYSDEC and the NYSDOH to collect a new sample in the heating season of 2020/2021 once additional passive mitigation measures were implemented (e.g., weather stripping and HVAC adjustments).

3 Building 3 Air Sampling

This report summarizes the programmatic sampling activities conducted in 2021, including the routine Building 3 sampling event on January 14, the subsequent installation and efficacy testing as well as supplemental weekly and system sampling, and the results of the ATS post-startup confirmation sampling event conducted on June 24 within the occupied boiler control room.

3.1 Pre-Sampling Chemical Inventory

Prior to the January and June 2021 sampling events, a detailed building review and reconnaissance was conducted to: (1) confirm and document specific sample locations (i.e., measured from outside walls); (2) obtain additional information on building layout, uses, and HVAC system configuration and operation; and (3) identify and remove chemical products from the building, if feasible. Chemical products that could not be removed from buildings 48 hours in advance of the sampling were documented. A photoionization detector (PID) equipped with an 11.7 electron volt (eV) lamp was used to screen for the presence of detectable vapor-phase chemicals during the reconnaissance. In addition, during the supplemental ATS weekly testing activities noted above, a chemical inventory was conducted in the vicinity (basement level) of the ATS.

Based on the January chemical inventory presented in **Attachment A**, one product was identified as containing TCE during the January inventory; a used 18-ounce bottle of *Nu-Calgon's*® Cal-Spray branded *nu-blast*® condenser/coil cleaner was identified per the Safety Data Sheet (SDS) to contain approximately 90-98% TCE by weight. This product was identified within a flammable storage cabinet located in Building 2 (which is adjacent to Building 3). Depending on when this product was last used, the residual concentration on the surface(s) it was applied to could continue to off-gas into the indoor air for an undetermined amount of time. This potential confounding source coincides with the low-level detection of TCE in the indoor air results. Naphthalene was also confirmed to be an ingredient in at least one product (i.e., Chesterton 730 Spragrip®) identified in Building 2. Both products had been removed from the building prior to the ATS start-up and subsequent sampling events.

As these products are part of BMS's chemical inventory and BMS has a Hazardous Materials Communication Program for workers withing Building 2 and 3, any exposures associated with this product would fall under the Occupational Safety and Health Administration (OSHA) worker exposure criteria of 50 parts per million (ppm) for 8-hour Threshold Limit Values (TLV). Therefore, further mitigation measures associated with the potential for SVI would not be required. Nevertheless, BMS will continue to operate the ATS to provide an additional level of protection for the workers within the boiler control room.

3.2 Air Sampling

Building 3 air sampling events were conducted in January, April, May and June 2021. With the January and June sampling events identified as the ATS pre- and post-startup (i.e., confirmation) sampling events for the ATS, respectively. As discussed in Section 5.3 below, in addition to the confirmatory air sampling events, supplemental weekly system testing samples and ATS influent/effluent air samples were collected during the first month of ATS operation.

3.2.1 Sample Types and Locations

As with previous Building 3 VI sampling events, the January and June 2021 sampling events each included collection of one indoor air (with duplicate) sample in the boiler control room and one outdoor ambient air sample. The 6-liter Summa canisters were collected over an 8-hour period at the same sampling locations where indoor and outdoor ambient air samples were collected historically and had been approved in the field with NYSDEC staff. The weekly ATS testing samples were also collected over and 8-hour sample duration from location IA-3 directly above the ATS. The influent and effluent samples were collected as a grab/instantaneous system sample from the corresponding side of the ATS from a sample port installed into the temporary ducting extensions. The sampling locations are shown on **Figure 3**.

3.2.2 Sampling Procedures and Analysis

Pre- and post-startup samples and associated quality assurance/quality control (QA/QC) samples were collected in accordance with the NYSDEC-approved *Field Sampling and Analysis Plan* (FSAP; OBG 2013b) and *Quality Assurance Project Plan* (QAPP; OBG 2013c) and the NYSDEC Division of Environmental Remediation's *DER-10 Technical Guidance for Site Investigation and Remediation* (DER-10; NYSDEC 2010). All work was performed under the responsible charge of a qualified environmental professional as defined in DER-10.

Indoor and ambient air samples were collected concurrently during the January 14 and June 24, 2021, sampling events, respectively. All air samples were collected in individually certified clean 6-liter (L) Summa® canisters with the indoor air and ambient air samples placed approximately three to five feet above the ground surface (i.e., at approximate breathing zone height) and collected over an 8-hour period, consistent with previous sampling events. No potential equipment quality control issues were identified during indoor or ambient air sampling.

All sample canisters were provided by Eurofins Air Toxics Laboratories Environmental, LLC, of Folsom, California. After sampling, the canisters were returned to the laboratory by overnight courier, under chain-of-custody, for analysis of VOC in accordance with USEPA Method TO-15. Additionally, as feasible, site-specific VOC not included in the laboratory's standard TO-15 compound list were reported by the laboratory as tentatively identified compounds (TIC).

4 Pre- and Post- ATS Air Sample Results

4.1 Applicable Screening Criteria

To evaluate the potential need for additional evaluation or mitigation, indoor and ambient air analytical data were compared to the applicable screening values. In accordance with DER-10 (Section 3.14(c)(4)), the analytical data was also compared to residential screening values. The current and reasonably anticipated future use of BMS buildings is commercial/industrial.

- For constituents of concern (COCs) included in the NYSDOH May 2017 decision matrices, the indoor air and soil vapor guidance values provided in the matrices are used to identify whether no further action, additional monitoring, or mitigation is needed, regardless of commercial or residential building occupancy.
- The Commercial Indoor Air Screening Level is the USEPA Building Assessment Survey and Evaluation (BASE) Study 90th percentile value, provided in Appendix C of the NYSDOH Guidance document, except for

- COCs included in the NYSDOH decision matrices. If a BASE value is not available, the USEPA Industrial Air Regional Screening Level (RSL) (USEPA 2021) using the lower of a target cancer incremental risk of 1x10⁻⁶ or a target hazard quotient of 1 is used, when available.
- NYSDOH Guidance document appendix, the Residential Indoor Air Screening Level is the NYSDOH Fuel Oil
 Study Upper Fence value, except for COCs included in the NYSDOH decision matrices. If a Fuel Oil Study
 Upper Fence value is not available, the USEPA Residential Air RSL value using the lower of a target cancer
 incremental risk of 1x10⁻⁶ or a target hazard quotient of 1 is used, when available. If a USEPA Residential Air
 RSL value is not available, the USEPA Industrial Air RSL is used, when available.

4.2 Comparison to Screening Criteria

As shown on **Table 1**, during the January 14, 2021 pre-startup sampling event, 32 VOC compounds were detected in either the parent or duplicate air samples; only four compounds exceeded either the commercial or residential screening values. During the June 24, 2021 post-startup (i.e., confirmation) sampling event, 24 compounds were detected, and only two exceeded screening values. A discussion of the compounds that exceeded screening values is presented below.

- Carbon tetrachloride was detected in both sampling rounds with almost identical concentrations between the ambient air sample and the indoor air sample. In both the January and June 2021 sampling events, carbon tetrachloride was detected at concentrations of 0.44 μg/m³ and 0.32 μg/m³ from the IA-3 parent sample and at 0.42 μg/m³ and 0.33 μg/m³ in the duplicate sample, which is above the NYSDOH decision matrix guidance value for monitoring of 0.2 μg/m³. However, outside ambient levels were almost identical at reported concentrations of 0.46 μg/m³ and 0.40 μg/m³, respectively. As such, the exceedances of the screening criteria in indoor air may be considered directly related to background influence from ambient air integration through the HVAC system.
- For just the January 2021 sampling event, ethanol was detected at 580 µg/m³ from the IA-3 parent sample and at 550 µg/m³ in the duplicate sample, which was below the Commercial Indoor Air Screening level of 1,300 µg/m³, but above the Residential Indoor Air Screening level of 210 µg/m³. Various concentrations of ethanol had been detected in four of the previous sampling events; however, not at such elevated levels. Since ethanol was not identified in groundwater directly upgradient from Building 3, the elevated detection was likely the result of ethanol-based hand/surface sanitizer being used by workers within the boiler control room. Ethanol did not exceed screening values in the June 2021 sampling event.
- Naphthalene was detected during both the January and June 2021 sampling events with estimated concentrations of 0.22 μg/m³ and 0.43 μg/m³ from the IA-3 parent sample and 0.26 μg/m³ and 0.42 μg/m³ in the duplicate sample, respectively. All samples were below the Commercial Indoor Air Screening level of 5.1 μg/m³ but above the Residential Indoor Air Screening value of 0.083 μg/m³. Varying concentrations of naphthalene were detected in each of the three previous sampling events. Since naphthalene was not identified in groundwater directly upgradient from Building 3, it is likely that the detection identified in the indoor air sample was the result of a transient source. Naphthalene is a known ingredient in many paints, oils, and fuels, and can off-gas from a worker's clothing or contaminated materials long after contact.
- TCE was detected at a concentration of 0.54 μg/m³ from the IA-3 parent sample and 0.52 μg/m³ from the
 duplicate sample during the January 2021 pre-startup sampling event. This TCE concentration was lower
 than any of the previous indoor air samples collected in Building 3, but greater than the NYSDOH decision
 matrix guidance value for monitoring (in the absence of companion soil vapor concentrations) of 0.2 μg/m³.

Following installation of the ATS, the indoor air TCE concentration from the June 2021 post-startup confirmation sampling event was lower than the NYSDOH decision matrix guidance value for monitoring; 0.16 μ g/m³ from the IA-3 parent sample and 0.15 μ g/m³ from the duplicate sample.

4.2.1 Data Usability

Upon completion of analysis, the laboratory data packages were validated by Arcadis. Laboratory analytical reports and corresponding Data Usability Summary Report (DUSR) for indoor air and ambient air samples are presented in **Attachment C** for each of the sampling events. No sample analytical results were rejected.

5 Mitigation Measures

5.1 Pre- ATS Installation Measures

In January 2021, facility HVAC settings were adjusted to allow the maximum fresh air mixing into the building. Measurements from BMS's HVAC contractor Jones Lang LaSalle Incorporated (JLL) were collected from each of the supply and return ducts and indicated a total of 2,659 cfm from the supply ducts and a total of 821 cfm from the return ducts; this indicated 1,838 cfm supplied from fresh air make-up. These results far exceed the United States Environmental Protection Agency (USEPA) guidance as referenced by American Society of Heating, Refrigerating and Air-Conditioning Engineers (ASHRAE) Standard 62.1-2001 to provide a minimum of 15 cfm of outdoor air per building occupant, considering building occupancy in Building 3 maintains no more than 10 people in each day, in the 1,838 cfm far exceeds the 150 cfm requirement if 10 employees were to be working in the building.

Following HVAC setting adjustments, differential pressure measurements were collected in January 2021 using stationary OmniGuard 4® differential pressure monitors to assess pressure differences between the Building 3 boiler control room indoor air and the following locations:

- Building 2 (directly connected to Building 3) indoor air;
- Inside a floor drain in the boiler control room, and
- Outdoor ambient air.

It should be noted that the monitored floor drain is typically fit with a one-way valve that was installed as part of initial mitigation measures implemented in 2018. The one-way valve was temporarily removed from the floor drain during differential pressure measurements.

As shown on **Attachment B**, the January results indicate that a positive pressure continues to exist within the boiler control room relative to ambient conditions and the adjoining Building 2. While a slight negative pressure (-0.002 inches of water column [iwc] to -0.005 iwc) was measured in the boiler control room indoor air relative to the floor drain feature, the pressure differential represented an improvement from measurements (-0.006 iwc in January 2019 taken prior to implementation of HVAC settings adjustments. The one-way valve was reinstalled in the floor drain following differential pressure monitoring mitigating potential VI.

In addition, during the spring of 2021, JLL (per BMS direction) installed micromanometer sensors in Building 3 to provide daily data records of differential pressure readings between the Building 3 basement boiler control room indoor air, and both the adjoining Building 2 indoor air and outside ambient air. This inital data set following data

recording set up is also presented in **Attachment B** and will be reported to the NYSDEC/NYSDOH within the monthly progress report submittals. The results indicate continuous positive pressure exist within the boiler control room relative to the rest of Building 3 (second floor) with some variability in pressure as it relates to ambient conditions and the adjoining Building 2.

5.2 ATS Installation Measures

As noted in Section 4 above, while a downward trend in indoor air TCE concentrations was demonstrated based on sampling conducted in January 2021 following mitigation measures installed in previous years, the detected TCE concentrations remained above the NYSDOH guidance value for monitoring of 0.2 µg/m³ for indoor air (NYSDOH, 2006). As a result, as shown on **Figure 3**, BMS opted to conservatively install an ATS in the boiler control room boiler control room as a supplemental mitigation measure to further reduce TCE concentrations in indoor air. The ATS consists of one variable-speed air treatment unit (Amaircare AirWash® MultiPRO) that includes an activated carbon canister for adsorption of VOC present in indoor air, as well as several particulate-filtering elements. The ATS was installed on April 22, 2021, and is intended to operate continuously. Based on post-startup field measurements, the variable speed range for the ATS (with the carbon canister) corresponds to a flow rate range of approximately 130 cubic feet per minute (cfm) to 300 cfm. The ATS has been operating at the low flow rate setting (130 cfm). An operation, maintenance, and monitoring (OMM) plan detailing the OMM procedures to be conducted for the ATS was submitted to the NYSDEC/NYSDOH on July 6, 2021.

Over the initial month of ATS operation, supplemental indoor air sampling was conducted and included the following:

- Eight-hour Indoor air samples via 6-liter Summa[®] canisters were collected in the boiler control room on a
 weekly basis to evaluate if changes in ATS operation (e.g., carbon breakthrough, speed adjustment) were
 warranted.
- An ATS influent and effluent air samples were also collected during the final weekly supplemental indoor air sampling event (i.e., about 1-month post-startup) to assess the removal efficiency of the carbon treatment.
 Temporary ductwork was installed on the ATS outlet to reduce potential dilution of the effluent sample.

Additional activities related to mitigation were conducted in January 2021 including the addition of weather stripping to the boiler control room doorways and balancing of the HVAC system to provide a more positive pressure environment.

5.3 Supplemental ATS Efficiently Air Sample Results

5.3.1 Weekly Indoor Air Samples

As shown in **Table 1**, weekly samples collected during the first month of ATS operation indicate a gradual reduction of TCE concentrations in indoor air. TCE concentrations ranged from an estimated concentration of 0.053 J μ g/m³ (1-month post-startup) to 0.28 μ g/m³ (1-week post-startup). As mentioned previously, supplemental indoor air sampling was conducted to confirm the system was operating at an appropriate setting (e.g., flow rate).

5.3.2 Influent/Effluent Air Samples

TCE was detected at an estimated concentration of 0.058 J ug/m3 in the ATS influent sample collected on May 20, 2021, approximately one month after the ATS began operating. As expected, assuming relatively consistent indoor air mixing in the boiler control room, this concentration was similar to the TCE concentration detected in the indoor air sample collected the same day $(0.053 \text{ J } \mu\text{g/m}^3)$. TCE was not detected in the ATS effluent sample. While using these data to quantify the carbon's removal efficiency may not be feasible (due to the relatively low influent concentrations), the influent/effluent sample results do support that the adsorption capacity of the carbon media was not spent at the time of the sampling.

6 Conclusions

Based on the results of indoor air sampling, the ATS has been effective at reducing indoor air concentrations of TCE in the boiler control room to below the NYSDOH decision matrix guidance value for monitoring of 0.2 µg/m³. The results also support that the low flow rate setting (approximately 130 cfm based on field measurements) is sufficient for achieving system performance criteria. Additionally, based on the June 2021 (2-months post-startup) and supplemental air sampling (i.e., weekly post-startup indoor air samples and influent/effluent ATS air samples) results, it appears that the carbon will maintain sufficient VOC removal efficiency beyond a period of 30 or 60 days (operating at the specified flow rate setting).

7 Future Activities

7.1 System Operation, Maintenance, and Monitoring

As detailed in the July 2021 Building 3 ATS OMM Plan, the first year of operation, indoor air sampling will be conducted monthly to support assessing the future carbon filter changeout schedule. Following the first year of operation, the indoor air sampling schedule will be reevaluated but is anticipated to change to biennial indoor air sampling during the heating season for 4 years followed by once every 4 years pending no increase in TCE levels have been identified and the ATS is meeting mitigation metrics.

During OMM Plan implementation, each ATS sampling event will consist of collecting one indoor air, one duplicate, and one ambient air sample using a passive sampler (e.g., Radiello 130/145 or similar). The use of an USEPA-approved sorbent-diffusive passive sampler is appropriate for this location due to the continuous air movement within the room and the long duration worker occupancy. The samplers provide a longer sample collection period (>24 hrs) which in turn will provide a better dataset for assessing long-term system functionality and the potential for human health exposures (USEPA 2014). The samplers will be placed approximately three to five feet above the floor or ground surface (i.e., at approximate breathing zone height) and collected over a 7-day period. Sample locations are shown on **Figure 3**. The ATS will be operating in the boiler control room with entry doors kept closed to the extent feasible for the duration of the sampling period. BMS will provide 7 days of notice to NYSDEC and NYSDOH in advance of indoor air sampling events.

Samples will be provided to an Environmental Laboratory Approval Program (ELAP)-certified laboratory for analysis of TCE in accordance with USEPA Method TO-17. One duplicate indoor air sample will be collected each day that indoor air sampling is performed. As applicable, all scheduled samples will be collected in accordance with current addendums to the NYSDEC-approved FSAP (OBG 2013a) and QAPP (OBG 2013b) and the NYSDEC Division of Environmental Remediation DER-10 (NYSDEC 2010).

For the first year of operation, the ATS carbon filter and particulate filters will be replaced on a quarterly basis. Following the first year of operation, the carbon filter replacement schedule will be reevaluated based on system performance and indoor air laboratory analytical results.

The effectiveness of the system will be evaluated based on indoor air sample results. TCE concentrations in indoor air will be compared to the NYSDOH guidance values of $0.2 \,\mu\text{g/m}^3$ for monitoring and $1.0 \,\mu\text{g/m}^3$ for mitigation (NYSDOH 2006). If over time indoor air concentrations fall below the TCE guidance value for monitoring ($0.2 \,\mu\text{g/m}^3$), modifications to the system will be considered and may include:

- Reducing indoor air sampling frequency
- Reducing carbon filter changeout frequency.

BMS will notify NYSDEC and NYSDOH in advance of implementing system optimization measures or modifications.

7.2 Reporting

Following the first year of ATS operation, a brief letter report will be provided to NYSDEC and NYSDOH after each boiler control room indoor air sampling event. The report will be submitted within 60 days of data validation and will include a summary of indoor air analytical data, OMM activities completed during the reporting period, and a summary of planned future OMM activities. In following years, BMS will submit an annual OMM report that includes the weekly and monthly checks of the system and the dates of completed maintenance. If sampling occurs during the year, a summary of the analytical data will also be submitted in the annual report.

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Tables



Building 3 VI and Air Treatment System Assessment Report Site #C734138: BMS Syracuse North Campus Restoration Area

		Indoor Air S	creening Value*	Building	a 3 - Mar. 2016	Building 3	- Dec. 2016	Building	3 - Jan. 2018	Building	3 - Jan. 2019	Building	3 - Jan. 2021		Build	ing 3 - ATS Te	stina April-M	lav 2021		Building 3 - J	June 2021
		maoor 7tir C		Ambient	Indoor	Ambient	Indoor	Ambient	Indoor	Ambient	Indoor	Ambient	Indoor	Indoor	Indoor	Indoor	Indoor	Indoor	Indoor	Indoor	Ambient
	CAS	Commercial (Exceedance		AA-3	IA-3	AA-3	IA-3	AA-3	IA-3	AMB-3	IA-3/DUP	AMB-011421	IA-3	IA-3	IA-3	IA-3	IA-3	INELLIENT-1	EFFLUENT-1	IA-3	AA-3
Date Collected:	Number	Gray Shaded		03/02/16	03/02/16	12/08/16	12/08/16	01/17/18	01/17/18	01/24/19	01/24/19	01/14/21	01/14/21	04/29/21	05/06/21	05/13/21	05/20/21	05/20/21	05/20/21	06/24/21	06/24/21
Volatile Organics - TO-15 (ug/m³)																					
1,1,1-Trichloroethane* 1,1,2,2-Tetrachloroethane	71-55-6 79-34-5	0.21	3 0.4	0.017 J 0.18 U	0.045 J [0.049 J] 0.20 U [0.20 U]	0.019 J 0.17 U	0.13 J 0.19 U	0.13 U 0.16 U	0.054 J [0.051 J] 0.19 U [0.19 U]	0.14 U 0.18 U	0.039 J [0.039 J] 0.18 U [0.19 U]	0.15 U 0.19 U	0.032 J [0.032 J]	0.16 U 0.20 U	0.72 U 0.91 U	1.6 U 2.0 U	0.014 J 0.19 U	0.014 J 0.16 U	0.0068 J	0.013 J [0.015 J]	0.0082 J 0.21 U
1,1,2-Trichloro-1,2,2-trifluoroethane	76-13-1	3.5	2.5	0.40 J	0.48 J [0.47 J]	0.56 J	0.60 J	0.52 J	0.57 J [0.55 J]	0.75 J	0.74 J [0.66 J]		0.10 0 [0.20 0]	0.200	0.010	2.00	0.10 0	0.100	0.100	0.10 0 [0.10 0]	0.210
(Freon 113)					1				1 1			0.43 J	0.52 J [0.47 J]	1.1 U	5.1 U	11 U	0.36 J	0.41 J	0.26 J	0.42 J [0.38 J]	0.46 J
1,1,2-Trichloroethane	79-00-5 75-34-3	0.7	0.4	0.14 U 0.10 U	0.16 U [0.16 U] 0.022 J [0.022 J]	0.14 U 0.10 U	0.15 U 0.052 J	0.13 U 0.097 U	0.15 U [0.15 U] 0.035 J [0.036 J]	0.14 U 0.11 U	0.15 U [0.15 U] 0.11 U [0.11 U]	0.15 U 0.11 U	0.15 U [0.16 U] 0.11 U [0.12 U]	0.16 U 0.12 U	0.72 U 0.54 U	1.6 U 1.2 U	0.15 U 0.11 U	0.13 U 0.094 U	0.15 U 0.11 U	0.15 U [0.15 U] 0.11 U [0.11 U]	0.16 U 0.12 U
1,1-Dichloroethene*	75-35-4		0.2	0.051 U	0.059 U [0.058 U]	0.050 U	0.063	0.048 U	0.066 [0.064]	0.053 U	0.053 U [0.054 U]	0.056 U	0.056 U [0.058 U]	0.057 U	0.26 U	0.57 U	0.056 U	0.046 U	0.053 U	0.056 U [0.055 U]	0.060 U
1,2,4-Trichlorobenzene 1,2,4-Trimethylbenzene	120-82-1 95-63-6	9.5	0.5 9.8	4.7 U 0.14 J	5.5 U [5.4 U] 0.16 J [0.17 J]	4.7 U 0.62 U	5.2 U 0.18 J	4.4 UJ 0.20 J	5.2 UJ [5.1 UJ] 0.34 J [0.37 J]	4.9 U 0.65 U	5.0 U [5.1 U] 10 [10]	5.2 U 0.15 J	5.3 U [5.4 U] 0.16 J [0.17 J]	5.4 U 0.71 U	25 U 3.3 U	7.0 U	5.3 U 0.19 J	4.3 U 0.20 J	5.0 U 0.66 U	5.3 U [5.2 U] 0.70 U [0.68 U]	5.6 U 0.74 U
1,2-Dibromoethane (EDB)	106-93-4	1.5	0.4	0.20 U	0.23 UB [0.22 UB]	0.19 UB	0.22 UB	0.18 U	0.21 U [0.21 U]	0.20 U	0.20 U [0.024 J]	0.22 U	0.027 J [0.22 U]	0.22 U	1.0 U	2.2 U	0.024 J	0.021 J	0.20 U	0.22 U [0.21 U]	0.23 U
1,2-Dichloro-1,1,2,2-tetrafluoroethane	76-14-2	6.8	0.4	0.10 J	0.10 J [0.11 J]	0.11 J	0.11 J	0.097 J	0.11 J [0.10 J]	0.12 J	0.11 J [0.15 J]	0.10 J	0.10 J [0.098 J]	0.11 J 0.87 U	0.93 U	2.0 U	0.12 J	0.12 J	0.15 J	0.092 J [0.095 J]	0.10 J
1,2-Dichlorobenzene 1,2-Dichloroethane	95-50-1 107-06-2	0.9	0.5	0.77 U 0.10 UB	0.89 U [0.88 U] 0.12 UB [0.12 UB]	0.76 U 0.10 UB	0.85 U 0.11 UB	0.72 U 0.056 J	0.84 U [0.82 U] 0.054 J [0.058 J]	0.80 U 0.088 J	0.80 U [0.82 U] 0.086 J [0.092 J]	0.84 U 0.10 J	0.85 U [0.88 U] 0.069 J [0.068 J]	0.051 J	4.0 U 0.54 U	8.6 U 1.2 U	0.85 U 0.057 J	0.70 U 0.087 J	0.80 U 0.037 J	0.85 U [0.84 U] 0.042 J [0.043 J]	0.91 U 0.049 J
1,2-Dichloropropane	78-87-5	1.6	0.4	0.59 U	0.68 U [0.68 U]	0.58 U	0.65 U	0.55 U	0.64 U [0.63 U]	0.61 U	0.62 U [0.63 U]	0.65 U	0.66 U [0.67 U]	0.67 U	3.1 U	6.6 U	0.66 U	0.26 J	0.62 U	0.66 U [0.64 U]	0.70 U
1,3,5-Trimethylbenzene 1,3-Butadiene	108-67-8 106-99-0	3.7	3.9 0.094	0.63 U	0.73 U [0.72 U]	0.62 U	0.69 U	0.59 U 0.26 U	0.68 U [0.67 U]	0.65 U 0.29 U	1.9 [2.0]	0.69 U 0.31 U	0.70 U [0.72 U] 0.31 U [0.32 U]	0.71 U 0.32 U	3.3 U 1.5 U	7.0 U 3.2 U	0.70 U 0.31 U	0.57 U 0.26 U	0.66 U 0.30 U	0.70 U [0.68 U]	0.74 U 0.33 U
1,3-Dichlorobenzene	541-73-1	2.4	0.094	0.28 U 0.77 U	0.33 U [0.099 J] 0.89 U [0.88 U]	0.28 U 0.76 U	0.31 U 0.85 U	0.26 U	0.31 <i>U</i> [0.30 <i>U</i>] 0.84 U [0.82 U]	0.29 U	0.30 U [0.30 U] 0.80 U [0.82 U]	0.84 U	0.85 U [0.88 U]	0.87 U	4.0 U	8.6 U	0.85 U	0.20 U	0.80 U	0.85 U [0.84 U]	0.91 U
1,4-Dichlorobenzene	106-46-7	5.5	1.2	0.15 UB	0.18 UB [0.18 UB]	0.15 UB	0.17 UB	0.14 U	0.17 UB [0.16 U]	0.16 UJ	0.16 UJ [0.16 UJ]		0.17 U [0.18 U]	0.17 U	0.80 U	1.7 U	0.17 U	0.14 U	0.16 U	0.17 U [0.17 U]	0.18 U
1,4-Dioxane 2,2,4-Trimethylpentane	123-91-1 540-84-1	2.5	0.56	0.46 U 3.0 U	0.53 U [0.53 U] 3.4 U [3.4 U]	0.45 U 2.9 U	0.51 U 3.3 U	0.43 U 2.8 U	0.50 U [0.49 U] 3.2 U [3.2 U]	0.48 U 3.1 U	0.48 U [0.49 U] 3.1 U [3.2 U]	0.047 J 3.3 U	0.065 J [0.13 J]	0.36 J 3.4 U	2.4 U 16 U	5.2 U 33 U	0.38 J 3.3 U	0.14 J 2.7 U	0.093 J 3.1 U	0.14 J [0.19 J]	0.098 J 3.5 U
2-Hexanone	591-78-6	130	31	2.6 U	3.0 U [3.0 U]	0.42 J	0.83 J	2.4 U	2.8 U [2.8 U]	2.7 U	2.7 U [0.95 J]	2.9 U	2.9 U [3.0 U]	3.0 U	14 U	29 U	2.9 U	2.4 U	2.7 U	2.9 U [0.19 J]	3.1 U
3-Chloropropene	107-05-1	2	0.47	2.0 UJ	2.3 UJ [2.3 UJ]	2.0 U	2.2 U	1.9 U	2.2 U [2.1 U]	2.1 U	2.1 U [2.1 U]	4.4 UJ	4.4 UJ [4.6 UJ]	2.3 U	10 U	22 U	2.2 U	1.8 U	2.1 U	2.2 UJ [2.2 UJ]	2.4 UJ
4-Ethyltoluene Acetone	622-96-8 67-64-1	3.6 98.9	115	0.63 UB 3.6 UB	0.73 UB [0.72 UB] 9.9 [5.4]	0.62 U 6.8	0.69 U 16	0.18 J 4.5	0.40 J [0.38 J] 6.4 [6.0]	0.65 U 7.9 J	2.3 [2.5] 7.3 J [9.7 J]	0.15 J 4.1	0.14 J [0.14 J] 7.9 [7.6]	0.71 U 22	3.3 U 11 J	7.0 U 11 J	0.15 J 27	0.17 J 24	0.66 U	0.70 U [0.68 U]	9.9
Benzene	71-43-2	9.4	13	0.44	0.60 [0.59]	0.36	0.46	0.58	0.66 [0.66]	0.56	0.64 [0.64]	0.65	0.67 [0.63]	0.46	0.21 J	2.3 U	0.22 J	0.23	0.21 UB	0.11 J [0.12 J]	0.10 J
Benzyl chloride (a-chlorotoluene)	100-44-7	6.8	0.057	0.66 U	0.77 U [0.76 U]	0.65 U	0.73 U	0.62 U	0.72 U [0.71 U]	0.69 U	0.69 U [0.71 U]	0.72 U	0.74 U [0.76 U]	0.75 U 0.97 U	3.4 U 4.4 U	7.4 U	0.74 U	0.60 U	0.69 U	0.74 U [0.72 U]	0.78 U
Bromodichloromethane Bromoform	75-27-4 75-25-2	0.33	0.076 2.6	0.86 U 1.3 U	0.41 J [0.38 J] 1.5 U [1.5 U]	0.84 U 1.3 U	0.19 J 1.4 U	0.80 U 1.2 U	0.93 <i>U</i> [0.92 <i>U</i>] 1.4 U [1.4 U]	0.89 U 1.4 U	0.90 <i>U</i> [0.92 <i>U</i>] 1.4 U [1.4 U]	0.94 U 1.4 U	0.057 J [0.061 J]	1.5 U	6.9 U	9.6 U 15 U	0.95 U 1.5 U	0.78 U 1.2 U	0.90 U 1.4 U	0.95 U [0.93 U] 1.5 U [1.4 U]	1.0 U 1.6 U
Bromomethane	74-83-9	1.7	0.5	2.5 U	2.9 U [2.8 U]	2.4 U	2.7 U	2.3 U	2.7 U [2.7 U]	2.6 U	2.6 U [2.7 U]	2.7 U	2.8 U [2.8 U]	2.8 UJ	13 UJ	28 UJ	2.8 U	2.2 U	2.6 U	2.8 U [2.7 U]	2.9 U
Carbon disulfide Carbon tetrachloride*	75-15-0 56-23-5	4.2	0.2	2.0 UJ 0.42	2.3 UJ [2.3 UJ] 0.42 [0.41]	2.0 U 0.46	2.2 U 0.45	1.9 U 0.38	2.2 U [2.1 U] 0.39 [0.39]	2.1 U 0.47	2.0 J [2.1 U] 0.47 [0.48]	2.2 U 0.46	2.2 U [2.3 U] 0.44 [0.42]	2.2 U 0.33	10 U 0.28 J	22 U 1.8 U	0.30	1.4 J 0.32	0.092 J	2.2 U [0.65 J] 0.32 [0.33]	2.4 U 0.40
Chlorobenzene	108-90-7	0.9	0.4	0.42 0.59 U	0.68 U [0.68 U]	0.58 U	0.45 0.65 U	0.55 U	0.64 U [0.63 U]	0.47 0.61 U	0.62 U [0.63 U]	0.64 U	0.65 U [0.67 U]	0.67 U	3.1 U	6.6 U	0.65 U	0.53 U	0.62 U	0.65 U [0.64 U]	0.70 U
Chloroethane	75-00-3	1.1	0.4	0.17 UJ	0.20 UJ [0.19 UJ]	0.17 U	0.19 U	0.16 U	0.18 U [0.18 U]	0.041 J	0.18 U [0.059 J]	0.027 J	0.041 J [0.046 J]	0.10 J	0.88 U	1.9 U	0.10 J	0.051 J		0.041 J [0.026 J]	0.030 J
Chloroform Chloromethane	67-66-3 74-87-3	3.7	1.2	0.080 J 0.91	1.4 [1.4] 0.91 [0.90]	0.086 J 0.78	1.4 0.79	0.089 J 0.80 J	0.74 [0.74] 0.85 J [0.83 J]	0.077 J 0.90 J	0.34 [0.33] 0.84 J [0.97 J]	0.070 J 0.62 J	0.38 [0.38] 0.64 J [0.64 J]	0.15 1.1 J	0.15 J 0.87 J	1.4 U 1.0 J	0.24 0.98 J	0.36 0.93 J	0.29 0.86 J	0.14 [0.15] 0.74 J [0.74 J]	0.065 J 0.71 J
cis-1,2-Dichloroethene*	156-59-2		0.2	0.10 U	0.14 [0.12]	0.10 U	0.29	0.095 U	0.28 [0.30]	0.10 U	0.12 [0.11]	0.11 U	0.069 J [0.064 J]	0.060 J	0.53 U	1.1 U	0.042 J	0.039 J	0.072 J	0.021 J [0.022 J]	0.12 U
cis-1,3-Dichloropropene	10061-01-5		0.4	0.58 U	0.67 U [0.67 U]	0.57 U	0.64 U	0.54 U	0.63 U [0.62 U]	0.60 U	0.61 U [0.62 U]	0.64 U	0.64 U [0.66 U]	0.66 U	3.0 U	6.5 U	0.64 U	0.53 U	0.61 U	0.64 U [0.63 U]	0.68 U
Cyclohexane Dibromochloromethane	110-82-7 124-48-1	26,000	6.3	0.44 U 1.1 U	0.51 U [0.50 U] 1.3 U [1.2 U]	0.43 U 1.1 U	0.098 J 1.2 U	0.15 J 1.0 U	0.29 J [0.26 J] 1.2 U [1.2 U]	0.46 U 1.1 U	0.46 U [0.47 U] 1.1 U [1.2 U]	2.4 U 1.2 U	2.4 U [2.5 U] 1.2 U [1.2 U]	2.5 U 1.2 U	11 U 5.7 U	25 U 12 U	2.4 U 1.2 U	0.36 J	2.3 U 1.1 U	2.4 U [1.6 J]	0.21 J
Dichlorodifluoromethane (Freon 12)	75-71-8	16.5	10	2.0	2.2 [2.2]	2.1	2.1	1.9	2.0 [2.0]	2.3 J	2.2 J [2.4 J]	2.3 J	2.3 J [2.3 J]	2.1 J	2.0 J	2.3 J	2.4 J	2.3 J	2.2 J	2.2 J [2.2 J]	2.1 J
Ethanol	64-17-5 100-41-4	1,300 5.7	210	2.5	20 [19]	2.7 0.074 J	66 0.13	3.4 0.16	72 [68]	5.3 J	51 J [57 J]	3.4	580 EJ [550 EJ]	3,000 EJ	1,600 EJ	3,300 EJ	1,100 EJ	560 EJ	840 EJ	100 J [100 J]	1.9 J
Ethylbenzene Hexachlorobutadiene	87-68-3	6.8	6.4 0.5	0.21 6.8 U	0.19 [0.30] 7.9 U [7.8 U]	6.7 U	7.5 U	6.4 UJ	0.20 [0.21] 7.4 UJ [7.3 UJ]	0.21 7.1 U	2.1 [1.8] 7.1 U [7.3 U]	0.10 J 7.5 U	0.11 J [0.10 J]	0.12 J 7.7 U	35 U	1.2 U 76 U	0.12 7.6 U	0.19	0.12 UB	0.058 J [0.065 J]	0.044 J 8.0 U
Isopropyl alcohol (2-propanol)	67-63-0	250	210	0.42 J	8.9 J [1.5 J]	0.36 J	4.5	1.5 U	3.2 [2.9]	1.3 J	4.6 J [5.3 J]	1.7 UB	19 [18]	130	68	82	42	49	88	6.2 [6.4]	2.6
Isopropylbenzene (cumene) m,p-Xylene	98-82-8 108-38-3	1,800 22.2	0.8	0.63 U 0.35	0.73 U [0.72 U] 0.56 [0.58]	0.62 U 0.22 UB	0.69 U 0.30	0.59 U 0.51	0.32 J [0.28 J] 0.57 [0.55]	0.65 U 0.65	0.40 J [0.67 U] 10 [9.1]	0.69 U 0.29	0.053 J [0.057 J] 0.35 [0.30]	0.71 U 0.38	3.3 U 0.28 J	7.0 U 2.5 U	0.70 U 0.39	0.073 J 0.52	0.66 U 0.23 UB	0.70 U [0.68 U] 0.20 J [0.22 J]	0.74 U 0.15 J
Methyl ethyl ketone (MEK, 2-Butanone)	78-93-3	12	16	0.66 J	1.1 J [0.46 J]	1.8 U	4.3	0.71 J	0.62 J [0.71 J]	1.9 J	1.2 J [2.3]	0.50 J	0.94 J [1.3 J]	1.5 J	9.8 U	21 U	1.6 J	3.9	2.0 U	1.1 J [1.4 J]	1.0 J
4-methyl-2-pentanone (MIBK)	108-10-1	6	1.9	0.52 U	0.61 U [0.60 U]	0.19 J	0.30 J	0.49 U	0.57 U [0.56 U]	0.22 J	0.23 J [0.56 U]	0.57 U	0.58 U [0.60 U]	0.59 U	2.7 U	5.8 U	0.19 J	0.21 J	0.55 U	0.14 J [0.15 J]	0.62 U
Methyl tert-butyl ether Methylene chloride*	1634-04-4 75-09-2	11.5	3	0.46 U 0.30 J	0.53 U [0.53 U] 1.0 U [0.29 J]	0.45 U 0.88 UB	0.51 U 0.98 UB	0.013 J 0.33 J	0.011 J [0.013 J] 0.64 J [0.42 J]	0.48 U 0.82 J	0.48 U [0.49 U] 0.64 J [0.58 J]	0.50 U 0.97 U	0.51 U [0.53 U] 0.99 UB [1.0 U]	0.52 U 0.69 J	2.4 U 4.6 U	5.2 U 9.9 U	0.51 U 2.5 U	0.032 J 0.92 J	0.48 U 2.3 U	0.51 U [0.50 U]	0.54 U 1.0 U
Naphthalene	91-20-3	5.1	0.083	0.34 U	0.39 U [0.38 U]	0.33 UB	0.23 J	0.31 UB	0.24 J [0.36 UB]	0.35 UB	4.6 J [4.2 J]	0.37 U	0.22 J [0.26 J]	0.22 J	1.7 U	3.7 U	0.61	0.67	0.35 U	0.43 [0.42]	0.12 J
n-Heptane	142-82-5	1,800	18	0.52 U	0.61 U [0.60 U]	0.52 UB	0.58 UB	0.21 J	0.36 J [0.70]	0.54 U	0.82 [0.99]	0.23 J	0.29 J [0.27 J]	3.0 U	14 U	29 U	0.60 J	0.50 J	2.7 U	2.9 U [2.8 U]	3.1 U
n-Hexane n-Propylbenzene	110-54-3 103-65-1	10.2 4,400	1.5	0.45 U 0.63 U	0.52 U [0.52 U] 0.73 U [0.72 U]	0.44 UB 0.62 U	0.50 UB 0.69 U	0.46 0.59 U	0.66 [1.0] 0.68 U [0.12 J]	0.76 0.65 U	0.61 [0.71] 1.2 [1.2]	0.36 J	0.54 J [0.42 J]	2.6 U 0.71 U	12 U 3.3 U	25 U 7.0 U	0.18 J	0.45 J 0.57 U	2.4 U 0.66 U	2.5 U [2.4 U] 0.70 U [0.68 U]	2.7 U 0.74 U
o-Xylene	95-47-6	7.9	7.1	0.13	0.21 [0.22]	0.084 J	0.12	0.22	0.24 [0.23]	0.32	4.1 [3.8]	0.11 J	0.12 [0.12 J]	0.15	0.58 U	1.2 U	0.15	0.22	0.013 J	0.071 J [0.093 J]	0.052 J
Styrene Tetropheroethope (PCE)*	100-42-5	1.9	1.4	0.54 U	0.63 U [0.63 U]	0.072 J	0.13 J	0.51 U	0.59 U [0.58 U]	0.57 U	0.57 U [0.58 U]	0.027 J	0.045 J [0.053 J]	0.52 J	2.8 U	6.1 U	0.14 J	0.078 J	0.081 J	0.070 J [0.093 J]	
Tetrachloroethene (PCE)* Tetrahydrofuran	127-18-4 109-99-9	8,800	0.8	0.039 J 1.9 U	0.077 J [0.081 J] 2.2 U [0.57 J]	0.044 J 1.8 U	0.11 J 0.52 J	0.075 J 1.8 U	0.14 J [0.18 U] 2.0 U [2.0 U]	0.067 J 2.0 U	0.18 J [0.17 J] 2.0 U [2.0 U]	0.13 J 2.1 U	0.74 [0.26] 2.1 U [2.2 U]	0.16 J 0.93 J	0.90 U 9.8 U	1.9 U 21 U	0.20 0.57 J	0.18 0.40 J	0.015 J 0.36 J	0.082 J [0.060 J] 2.1 U [0.22 J]	0.048 J 2.2 U
Toluene	108-88-3	43	57	0.87	1.5 [1.6]	0.41	0.77	0.82	1.0 [0.90]	0.81	5.0 [4.5]	0.65	0.67 [0.65]	0.58	0.72 J	0.70 J	0.52	16	0.018 J	0.45 [0.51]	0.24 J
trans-1,2-Dichloroethene trans-1,3-Dichloropropene	156-60-5 10061-02-6	 5 1.3	0.25	0.51 UJ 0.58 U	0.024 J [0.027 J] 0.67 U [0.67 U]	0.50 U 0.57 U	0.56 UB 0.64 U	0.48 U 0.54 U	0.23 J [0.22 J] 0.63 U [0.62 U]	0.53 U 0.60 U	0.53 U [0.54 U] 0.61 U [0.62 U]	0.56 U 0.64 U	0.56 U [0.58 U] 0.64 U [0.66 U]	0.57 U 0.66 U	2.6 U 3.0 U	5.7 U 6.5 U	0.56 U 0.64 U	0.012 J 0.53 U	0.013 J	0.56 U [0.55 U] 0.64 U [0.63 U]	0.60 U 0.68 U
Trichloroethene (TCE)*	79-01-6		0.25	0.58 U	0.67 0 [0.67 0] 0.93 [0.93]	0.57 U 0.14 UB	1.6	0.54 U	0.63 0 [0.62 0] 0.77 [0.82]	0.033 J	0.59 [0.61]	0.023 J	0.54 [0.52]	0.88	0.26 J	0.51 UB	0.053 J	0.058 J	0.61 U	0.16 [0.15]	0.017 J
Trichlorofluoromethane (Freon 11)	75-69-4	18.1	12	1.0	1.2 [1.1]	1.2	1.2	1.0	1.2 [1.0]	1.4	1.2 [1.2]	1.2	1.4 [1.2]	1.3	1.1 J	8.0 U	1.4	1.4	1.9	1.1 [1.1]	1.2
Vinyl chloride*	75-01-4		0.2	0.033 U	0.038 U [0.038 U]	0.032 U	0.018 J	0.031 U	0.036 U [0.035 U]	0.034 U	0.034 U [0.035 U]	0.036 U	0.036 U [0.037 U]	0.037 U	0.17 U	0.36 U	0.036 U	0.030 U	0.034 U	0.036 U [0.036 U]	0.038 U
Volatile Organics-TIC (ppbv) Isobutanol	78-83-1	4,400	1,000	NR	1.8 JN [2 JN]	NR	NR	0.79 JN	1.6 JN [1.9 JN]	NR	0.93 JN [1.6 JN]	NR	NR	NR	NR	NR	NR	1.1 NJ	NR	NR	NR
Pentane	109-66-0	260	2.5	NR	NR	NR	NR	NR	NR	NR	0.89 JN [1.2 JN]	NR	1.6 [1.3]	NR	NR	NR	1.1 NJ	5.8 NJ	NR	1.6 NJ [1.6 NJ]	NR
Propylene	526-73-8			NR	NR	NR	NR	NR	NR	NR	NR	NR	NR	2.1 NJ	NR	NR	NR	NR	NR	NR	NR

See Notes on Page 2.

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https://arcadiso365.sharepoint.com/leams/BMS-ESyracuseBDAProgram/Shared Documents/General/BLD 3 SVI/2021 Bldg 3 VI-ATS Report/Bldg 3 VI-ASSessment Rpt_Table 1 & 2

Table 1

Building 3 Air Analytical Results



Building 3 VI and Air Treatment System Assessment Report

Site #C734138: BMS Syracuse North Campus Restoration Area

Notes:

- 1. Samples were collected by Arcadis and analyzed by Eurofins Air Toxics Laboratories Environmental, LLC of Folsom, CA.
- 2. * = constituent included in the NYSDOH May 2017 Soil Vapor / Indoor Air Matrices indoor air guidance values listed guidance may be used to identify whether to conduct additional monitoring or mitigation.
- 3. The Commercial Indoor Air Screening Level is the USEPA BASE Study 90th percentile value, when available (except
- 4. The Residential Indoor Air Screening Level is the NYSDOH Fuel Oil Study Upper Fence value (except for *
- 4. Non-numerical values in the "CAS Number" column are a surrogate identification because no actual CAS number is available.
- 5. Analytes detected in sample are shown in black font and analytes that are not detected are shown in gray font.
- 6. Field duplicate sample results are presented in brackets, [].
- 7. The data has been validated.
- 8. Designations:
 - a) Italic font = Sample MDL exceeds the constituent's lower indoor air screening value.
 - Black bold font = Result detected above method detection limit
 - Gray shading = Result detected above constituent's higher indoor air screening value.

9. Abbreviations:

-- = Screening Value not available based on inquiry described in Notes 3 and 4.

 μ g/m³ = Micrograms per cubic meter.

CAS = Chemical Abstracts Service.

AA = Ambient Air

IA = Indoor Air

MDL = Method detection limit.

NR = Not reported as a TIC.

ppbv = Parts per billion by volume

10. Qualifier Definitions:

- B = Analyte was detected in the blank and sample.
- E = Analyte exceeded calibration range.
- J = Estimated value. Result is greater than the MDL but less than the RL.
- N = The analysis indicates the presence of a compound which there is presumptive evidence to make a tentative identification.
- U = Analyte not detected above the method detection limit. The compound reporting limit is presented for reference.
- UB = Analyte considered non-detect at the listed value due to associated blank contamination.
- UJ = The analyte was not detected above the reported sample detection limit. However, the reported limit is approximate and may or may not represent the actual limit of detection.

Table 2
Upgradient Groundwater Results



Building 3 VI and Air Treatment System Assessment Report Site #C734138: BMS Syracuse North Campus Restoration Area

Location ID:		TOGS 1.1.1			BDA-1F					BDA-1WT		
Date Collected:		Groundwater	04/12/16	07/17/17	12/08/17	09/28/18	10/30/19	04/12/16	07/17/17	12/08/17	09/27/18	10/30/19
	CAS	Standard/Guidance	BDA-1F	BDA-1F	BDA-1F	BDA-1F	BDA-1F	BDA-1WT	BDA-1WT	BDA-1WT	BDA-1WT	BDA-1WT
Sample Name:	Number	Value	04122016	07172017	12082017	09282018	10302019	04122016	07172017	12082017	09272018	10302019
Volatile Organics - USEPA SW-846 Method 8260C												
1,1,1-Trichloroethane	71-55-6	5	1.0 U [1.0 U]	1.0 U	1.0 U	1.0 U [1.0 U]	1 U	1.0 U	1.0 U	1.0 U	1.0 U	1 U
1,1,2,2-Tetrachloroethane	79-34-5	5	1.0 U [1.0 U]	1.0 U	1.0 U	1.0 U [1.0 U]	1 U	1.0 U	1.0 U	1.0 U	1.0 U	1 U
1,1,2-Trichloro-1,2,2-trifluoroethane	76-13-1	5	10 U [10 U]	10 U	10 U	10 U [10 U]	10 U	10 U	10 U	10 U	10 U	10 U
1,1,2-Trichloroethane	79-00-5	1	1.0 U [1.0 U]	1.0 U	1.0 U	1.0 U [1.0 U]	1 U	1.0 U	1.0 U	1.0 U	1.0 U	1 U
1,1-Dichloroethane	75-34-3	5	2.0 [2.0]	1.0 U	1.0 U	1.0 U [1.0 U]	1 U	1.0 U	1.0 U	1.0 U	1.0 U	1 U
1,1-Dichloroethene	75-35-4	5	2.0 [1.0]	1.0 U	1.0 U	1.0 U [1.0 U]	1 U	1.0 U	1.0 U	1.0 U	1.0 U	1 U
1,2,3-Trichlorobenzene	87-61-6	5	5.0 U [5.0 U]	5.0 U	5.0 U	5.0 U [5.0 U]	5 U	5.0 U	5.0 U	5.0 U	5.0 U	5 U
1,2,4-Trichlorobenzene	120-82-1	5	5.0 U [5.0 U]	5.0 U	5.0 U	5.0 U [5.0 U]	5 U	5.0 U	5.0 U	5.0 U	5.0 U	5 U
1,2,4-Trimethylbenzene	95-63-6	5	5.0 U [5.0 U]	5.0 U	5.0 U	5.0 U [5.0 U]	5 U	5.0 U	5.0 U	5.0 U	5.0 U	5 U
1,2-Dibromo-3-chloropropane	96-12-8	0.04	5.0 U [5.0 U]	5.0 U	5.0 U	5.0 U [5.0 U]	5 U	5.0 U	5.0 U	5.0 U	5.0 U	5 U
1,2-Dibromoethane	106-93-4	0.0006	1.0 U [1.0 U]	1.0 U	1.0 U	1.0 U [1.0 U]	1 U	1.0 U	1.0 U	1.0 U	1.0 U	1 U
1,2-Dichlorobenzene	95-50-1	3	5.0 U [5.0 U]	5.0 U	5.0 U	5.0 U [5.0 U]	5 U	5.0 U	5.0 U	5.0 U	5.0 U	5 U
1,2-Dichloroethane	107-06-2	0.6	1.0 U [1.0 U]	1.0 U	1.0 U	1.0 U [1.0 U]	1 U	1.0 U	1.0 U	1.0 U	1.0 U	1 U
1,2-Dichloropropane	78-87-5	1	1.0 U [1.0 U]	1.0 U	1.0 U	1.0 U [1.0 U]	1 U	1.0 U	1.0 U	1.0 U	1.0 U	1 U
1,3,5-Trimethylbenzene	108-67-8	5	5.0 U [5.0 U]	5.0 U	5.0 U	5.0 U [5.0 U]	5 U	5.0 U	5.0 U	5.0 U	5.0 U	5 U
1,3-Dichlorobenzene	541-73-1	3	5.0 U [5.0 U]	5.0 U	5.0 U	5.0 U [5.0 U]	5 U	5.0 U	5.0 U	5.0 U	5.0 U	5 U
1,4-Dichlorobenzene	106-46-7	3	5.0 U [5.0 U]	5.0 U	5.0 U	5.0 U [5.0 U]	5 U	5.0 U	5.0 U	5.0 U	5.0 U	5 U
1,4-Dioxane	123-91-1		250 U [250 U]	250 U	250 U	250 U [250 U]	250 U	250 U	250 U	250 U	250 U	250 U
2-Hexanone	591-78-6	50	10 U [10 U]	10 U	10 U	10 U [10 U]	10 U	10 U	10 U	10 U	10 U	10 U
2-Nitropropane	79-46-9		10 U [10 U]	10 U	10 U	10 U [10 U]	10 U	10 U	10 U	10 U	10 U	10 U
Acetone	67-64-1	50	20 U [20 U]	20 U	20 U	2.0 J [2.0 J]	20 U	20 U	20 U	20 U	0.90 J	20 U
Acetonitrile	75-05-8		100 U [100 U]	100 U	100 U	100 U [100 U]	100 U	100 U	100 U	100 U	100 U	100 U
Benzene	71-43-2	1	1.0 U [1.0 U]	1.0 U	1.0 U	1.0 U [1.0 U]	1 U	1.0 U	1.0 U	1.0 U	1.0 U	1 U
Bromochloromethane	74-97-5	5	5.0 U [5.0 U]	5.0 U	5.0 U	5.0 U [5.0 U]	5 U	5.0 U	5.0 U	5.0 U	5.0 U	5 U
Bromodichloromethane	75-27-4	50	1.0 U [1.0 U]	1.0 U	1.0 U	1.0 U [1.0 U]	1 U	1.0 U	1.0 U	1.0 U	1.0 U	1 U
Bromoform	75-25-2	50	4.0 U [4.0 U]	4.0 U	4.0 U	4.0 U [4.0 U]	4 U	4.0 U	4.0 U	4.0 U	4.0 U	4 U
Bromomethane	74-83-9	5	1.0 U [1.0 U]	1.0 U	1.0 U	1.0 U [1.0 U]	1 U	1.0 U	1.0 U	1.0 U	1.0 U	1 U
Carbon disulfide	75-15-0	60	5.0 U [5.0 U]	5.0 U	5.0 U	5.0 U [5.0 U]	5 U	5.0 U	5.0 U	5.0 U	5.0 U	5 U
Carbon tetrachloride	56-23-5	5	1.0 U [1.0 U]	1.0 U	1.0 U	1.0 U [1.0 U]	1 U	1.0 U	1.0 U	1.0 U	1.0 U	1 U
Chlorobenzene	108-90-7	5	1.0 U [1.0 U]	1.0 U	1.0 U	1.0 U [1.0 U]	1 U	1.0 U	1.0 U	1.0 U	1.0 U	1 U
Chlorodifluoromethane	75-45-6	5	5.0 U [5.0 U]	5.0 U	5.0 UJ	5.0 U [5.0 U]	5 UJ	5.0 U	5.0 U	5.0 UJ	5.0 U	5 UJ
Chloroethane	75-00-3	5	0.90 J [0.90 J]	1.0 U	1.0 U	1.0 U [1.0 U]	1 UJ	1.0 U	1.0 U	1.0 U	1.0 U	1 UJ
Chloroform	67-66-3	7	1.0 U [1.0 U]	1.0 U	1.0 U	0.70 J [0.60 J]	0.2 J	1.0 U	1.0 U	1.0 U	1.0 U	1 U
Chloromethane	74-87-3	5	1.0 U [1.0 U]	1.0 U	1.0 U	1.0 U [1.0 U]	1 UJ	1.0 U	1.0 U	1.0 U	1.0 U	1 UJ
cis-1,2-Dichloroethene	156-59-2	5	2.0 [2.0]	1.0	0.70 J	0.70 J [0.70 J]		1.0 U	1.0 U	1.0 U	1.0 U	1 U
cis-1,3-Dichloropropene	10061-01-5		1.0 U [1.0 U]	1.0 U	1.0 U	1.0 U [1.0 U]	1 U	1.0 U	1.0 U	1.0 U	1.0 U	1 U
Cyclohexane	110-82-7		5.0 U [5.0 U]	5.0 U	5.0 U	5.0 U [5.0 U]	5 U	5.0 U	5.0 U	5.0 U	5.0 U	5 U
Cyclohexanone	108-94-1		100 U [100 U]	100 U	100 UJ	100 U [100 U]	100 U	100 U	100 U	100 UJ	100 U	100 U
Dibromochloromethane	124-48-1	50	1.0 U [1.0 U]	1.0 U	1.0 U	1.0 U [1.0 U]	1 U	1.0 U	1.0 U	1.0 U	1.0 U	1 U
Dichlorodifluoromethane	75-71-8	5	1.0 U [1.0 U]	1.0 U	1.0 U	1.0 U [1.0 U]	1 UJ	1.0 U	1.0 U	1.0 U	1.0 U	1 UJ
Ethyl actate	141-78-6		5.0 U [5.0 U]	5.0 U	5.0 U	5.0 U [5.0 U]	5 U	5.0 U	5.0 U	5.0 U 5.0 U	5.0 U	5 U
Ethyl ether	60-29-7 100-41-4		5.0 U [5.0 U]	5.0 U	5.0 U	5.0 U [5.0 U]	5 U	5.0 U	5.0 U		5.0 U	5 UJ 1 U
Ethylbenzene	78-83-1	5	1.0 U [1.0 U]	1.0 U	1.0 U	1.0 U [1.0 U]	1 U	1.0 U	1.0 U	1.0 U	1.0 U	
Isobutanol	98-82-8		250 U [250 U]	250 U	250 U	250 U [250 U]	250 U	250 U	250 U	250 U	250 U	250 U
Isopropylbenzene		5	5.0 U [5.0 U]	5.0 U	5.0 U	5.0 U [5.0 U]	5 U	5.0 U	5.0 U	5.0 U	5.0 U	5 U
m,p-Xylene	179601-23-1		1.0 U [1.0 U]	1.0 U	1.0 U	5.0 U [5.0 U]	5 U	1.0 U	1.0 U	1.0 U	5.0 U	5 U
Methyl actual ketana (MEK, 2 butanana)	79-20-9		5.0 U [5.0 U]	5.0 U	5.0 U	5.0 U [5.0 U]	5 U	5.0 U	5.0 U	5.0 U	5.0 U	5 U
Methyl ethyl ketone (MEK, 2-butanone)	78-93-3	50	10 U [10 U]	10 U	10 U	10 U [10 U]	10 U	10 U	10 U	10 U	10 U	10 U

See Notes on Page 2.

Table 2 Upgradient Groundwater Results



Building 3 VI and Air Treatment System Assessment Report Site #C734138: BMS Syracuse North Campus Restoration Area

Location ID:		TOGS 1.1.1			BDA-1F					BDA-1WT		
Date Collected:		Groundwater	04/12/16	07/17/17	12/08/17	09/28/18	10/30/19	04/12/16	07/17/17	12/08/17	09/27/18	10/30/19
	CAS	Standard/Guidance	BDA-1F	BDA-1F	BDA-1F	BDA-1F	BDA-1F	BDA-1WT	BDA-1WT	BDA-1WT	BDA-1WT	BDA-1WT
Sample Name:	Number	Value	04122016	07172017	12082017	09282018	10302019	04122016	07172017	12082017	09272018	10302019
/olatile Organics - USEPA SW-846 Method 8260C (μg/L) (cont'd)												
Methyl isobutyl ketone (MIBK, 4-methyl-2-pentanone)	108-10-1	† (50)	10 U [10 U]	10 U	10 U	10 U [10 U]	10 U					
Methyl tert-butyl ether	1634-04-4	10	1.0 U [1.0 U]	1.0 U	1.0 U	1.0 U [1.0 U]	1 U	1.0 U	1.0 U	1.0 U	1.0 U	1 U
Methylcyclohexane	108-87-2	†	5.0 U [5.0 U]	5.0 U	5.0 U	5.0 U [5.0 U]	5 U	5.0 U	5.0 U	5.0 U	5.0 U	5 U
Methylene chloride	75-09-2	5	4.0 U [4.0 U]	4.0 U	1.0	0.90 J [0.90 J]		4.0 U	4.0 U	1.0 U	1.0 U	1 U
Naphthalene	91-20-3	10	5.0 U [5.0 U]	5.0 U	5.0 U	5.0 U [5.0 U]	5 U	5.0 U	5.0 U	5.0 U	5.0 U	5 U
n-Butanol	71-36-3	50	250 U [250 U]	250 U	250 U	250 U [250 U]	250 U					
n-Butylbenzene	104-51-8	5	5.0 U [5.0 U]	5.0 U	5.0 U	5.0 U [5.0 U]	5 U	5.0 U	5.0 U	5.0 U	5.0 U	5 U
n-Heptane	142-82-5		5.0 U [5.0 U]	5.0 U	5.0 U	5.0 U [5.0 U]	5 U	5.0 U	5.0 U	5.0 U	5.0 UJ	5 U
n-Hexane	110-54-3		5.0 U [5.0 U]	5.0 U	5.0 U	5.0 U [5.0 U]	5 U	5.0 U	5.0 U	5.0 U	5.0 U	5 U
o-Xylene	95-47-6	5	1.0 U [1.0 U]	1.0 U	1.0 U	1.0 U [1.0 U]	1 U	1.0 U	1.0 U	1.0 U	1.0 U	1 U
p-Isopropyltoluene	99-87-6	5	5.0 U [5.0 U]	5.0 U	5.0 U	5.0 U [5.0 U]	5 U	5.0 U	5.0 U	5.0 U	5.0 U	5 U
sec-Butylbenzene	135-98-8	5	5.0 U [5.0 U]	5.0 U	5.0 U	5.0 U [5.0 U]	5 U	5.0 U	5.0 U	5.0 U	5.0 U	5 U
Styrene	100-42-5	5	5.0 U [5.0 U]	5.0 U	5.0 U	5.0 U [5.0 U]	5 U	5.0 U	5.0 U	5.0 U	5.0 U	5 U
tert-Butylbenzene	98-06-6	5	5.0 U [5.0 U]	5.0 U	5.0 U	5.0 U [5.0 U]	5 U	5.0 U	5.0 U	5.0 U	5.0 U	5 U
Tertiary butyl alcohol	75-65-0		20 U [20 U]	20 U	20 U	50 U [50 U]	50 U	20 U	20 U	20 U	50 U	50 U
Tetrachloroethene	127-18-4	5	1.0 U [1.0 U]	1.0 U	1.0 U	1.0 U [1.0 U]	1 U	1.0 U	1.0 U	1.0 U	1.0 U	1 U
Tetrahydrofuran	109-99-9	50	10 U [10 U]	10 U	10 U	10 U [10 U]	10 U					
Toluene	108-88-3	5	1.0 U [1.0 U]	1.0 U	1.0 U	1.0 U [1.0 U]	1 U	1.0 U	1.0 U	1.0 U	1.0 U	1 U
trans-1,2-Dichloroethene	156-60-5	5	1.0 U [1.0 U]	1.0 U	1.0 U	1.0 U [1.0 U]	1 U	1.0 U	1.0 U	1.0 U	1.0 U	1 U
trans-1,3-Dichloropropene	10061-02-6	0.4	1.0 U [1.0 U]	1.0 U	1.0 U	1.0 U [1.0 U]	1 U	1.0 U	1.0 U	1.0 U	1.0 U	1 U
Trichloroethene	79-01-6	5	14 [13]	11	5.0	7.0 [6.0]	2	1.0 U	1.0 U	1.0 U	1.0 U	1 U
Trichlorofluoromethane	75-69-4	5	1.0 U [1.0 U]	1.0 U	1.0 U	1.0 U [1.0 U]	1 UJ	1.0 U	1.0 U	1.0 U	1.0 U	1 UJ
Vinyl chloride	75-01-4	2	1.0 U [1.0 U]	1.0 U	1.0 U	1.0 U [1.0 U]	1 UJ	1.0 U	1.0 U	1.0 U	1.0 U	1 UJ
Xylenes (total)	1330-20-7	5	1.0 U [1.0 U]	1.0 U	1.0 U	5.0 U [5.0 U]	6 U	1.0 U	1.0 U	1.0 U	5.0 U	6 U

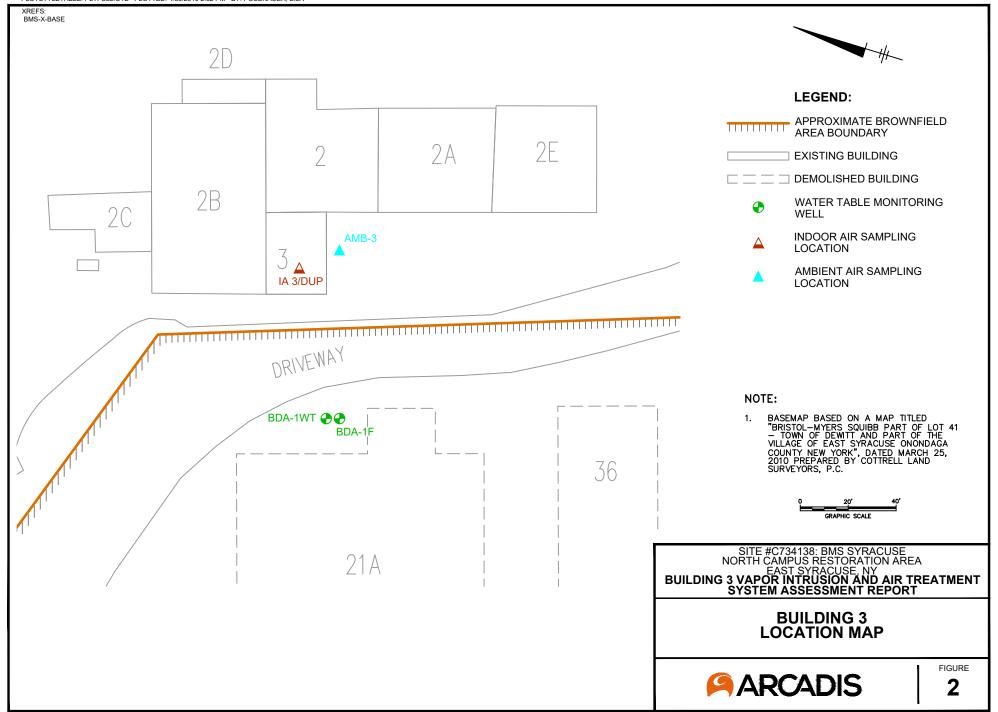
Notes:

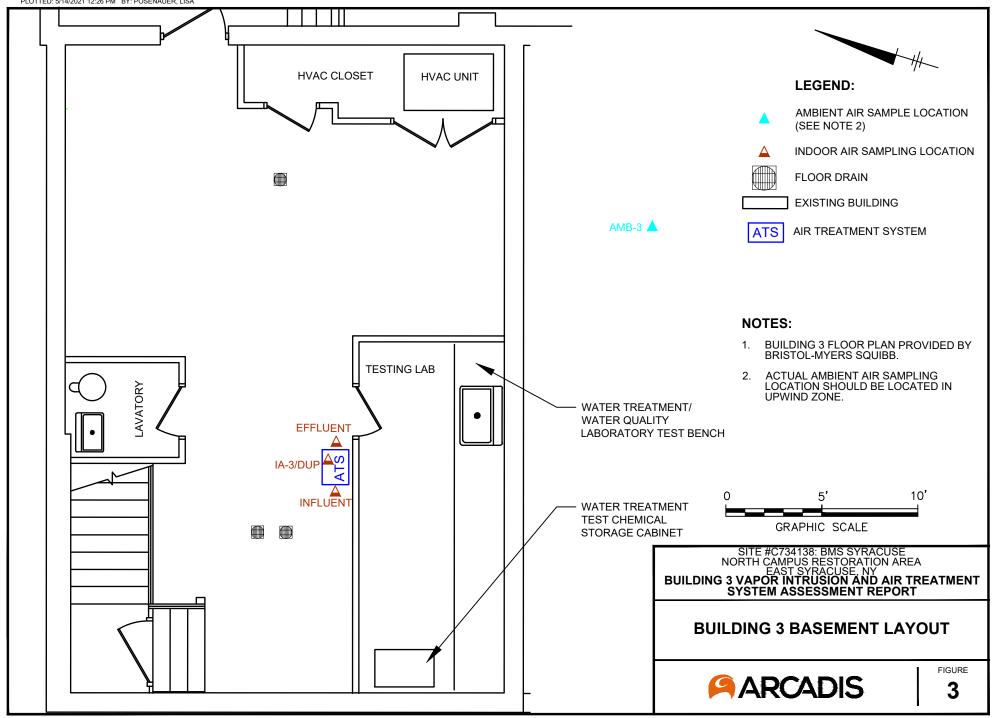
- 1. Results compared to NYSDEC Division of Water Technical and Operational Guidance Series (TOGS) 1.1.1, Ambient Water Quality Standards and Guidance Values.
- 2. The data has been validated.
- 3. Samples were analyzed by TestAmerica Laboratories of Buffalo/Amherst, New York or Eurofins Lancaster Laboratories Environmental, LLC, Lancaster, PA.
- 4. Analytes detected in sample are shown in black font and analytes that are not detected are shown in gray font.
- 5. Field duplicate sample results are presented in brackets, [].
- 6. The compounds 3-methylphenol and 4-methylphenol cannot be chromatographically separated, and therefore are reported as 3&4-methylphenol.
- 7. Data have been validated.
- 8. Designations:
 - a) Bold font with gray shading = Result detected at or above NYSDEC TOGS 1.1.1. or a default screening value for purposes of this report.
 - b) Italic font = Sample MDL exceeds NYSDEC TOGS 1.1.1 value.
 - c) † = Not determined to be a POC at this time based on review of TOGS Part 1(B)(2) Steps 1 through 3.
 - d) () = Value in parenthesis and gray font indicate a default screening value being used for purposes of this report.
- 9. Abbreviations:
 - -- = Substance not regulated by the POC Groundwater Standard (TOGS 1.1.1, Table 3).
 - μg/L = Micrograms per liter.
 - CAS = Chemical Abstracts Service.
 - NYSDEC = New York State Department of Environmental Conservation.
 - VOC = Volatile organic compound.
- 10. Qualifier Definitions:
 - J = Estimated value. Result is greater than the MDL but less than the RL.
 - U = Analyte not detected at listed reporting limit.

Figures

NEW YORK

360/akradiashaNa - BRISTOL-MYERS COMPANYProject FilesBDA Hydraulic Testing Program/2021/01-DWGB3OMM_F01_SLM.dwg LAYOUT: 1 SAVED: 5/13/2021 5/03 PM ACADVER: 24.0S (LMS TECH) PAGESETUP: C-PA-PDF PLOTSTYLETABLE: PLTFULL.CTB PLOTTED:





Appendix A

Structural Sampling Questionnaire and Building Inventory Forms/Photo Logs (See Attached CD for Inventory Photos)



Site Name: Bristol-Myers Squibb		Site Code:	Operable Unit:
Building Code: Industrial	Building Name:_	Building	3
Address: 6000 Thompson Rd			Apt/Suite No:
City: East Syracuse	State: NY	Zip: 13057	County: Onondaga
Contact Information			
Preparer's Name: Daniel Zuck			Phone No: 516-369-2741
Preparer's Affiliation: Arcadis of NY Inc.			Company Code:
Purpose of Investigation: Soil Vapor Investig	gation		Date of Inspection: 1/14/2021
Contact Name: Anne Locke		1111	Affiliation: MANAGER
Phone No: 315-432-2660 Alt. Phone	ne No:		Email: anne.locke@bms.com
Number of Occupants (total): 5-10 Number	r of Children: 0		
Cocupant Interviewed?	Owner Occup	pied?	☐ Owner Interviewed?
Owner Name (if different): Bristol-Myers Squi	ibb		Owner Phone: NA
Owner Mailing Address: PO Box 4755, Syracus	se NY 13221		
Building Details			
Bldg Type (Res/Com/Ind/Mixed): COMMERCIAL/N	MIXED		Bldg Size (S/M/L): MEDIUM
If Commercial or Industrial Facility, Select Operations: MULTI-USE BUILDING	If	Residential Sele	ect Structure Type:
Number of Floors: 2 Approx. Year Construc	tion:1910	🗷 Buildin	g Insulated? Attached Garage
Describe Overall Building 'Tightness' and Airflows(e.g., r	results of smoke tests):	u - 2 - 24 - 24 - 24 - 24 - 24 - 24 - 24
Building has multiple air handlers a infiltration can be detected at buil			doors. Ambient air
Foundation Description	turng decess io	cacions.	
Foundation Type: BASEMENT-PARTIAL	Fo	undation Deptl	n (bgs): 8 Unit: FEET
Foundation Floor Material: POURED CONCRETE	Fo	undation Floor	Thickness: 7
Foundation Wall Material: CONCRETE BLOCK	Fo	undation Wall 1	Thickness: Unit: INCHES
Floor penetrations? Describe Floor Penetrations:	Floor drains		30000
Wall penetrations? Describe Wall Penetrations:	Electrical co	onduits, pip	ping
Basement is: PARTIALLY FINISHE Basement is:	DRY	Sump:	s/Drains? Water In Sump?: YES
Describe Foundation Condition (cracks, seepage, etc.) :	Solid but cra	acked, Cove	ed by tile in spots.
Radon Mitigation System Installed?	VOC Mitigatio	n System Install	led? Mitigation System On
Heating/Cooling/Ventilation Systems			
Heating System: FORCED AIR	Heat Fuel Type:	GAS	▼ Central A/C Present?
Vented Appliances			
Vented Appliances Water Heater Fuel Type: ELECTRIC		thes Dryer Fuel	Type: NO CLOTHES DRYER



Structure Sampling Questionnaire and Building Inventory

New York State Department of Environmental Conservation

		PI	RODUCT INVEN	ITORY			
Building Nam	e: Building 3		1071200 1 (20 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1	e: Industrial	Date:	1/14/2023	1
Bldg Address	: 6000 Thompson Rd				Apt/Suite	No:	
Bldg City/Stat	t e/Zip: East Syracuse N	Y, 1305	7				
Make and Mo	del of PID: ppbRAE 2000			Date of	Calibration:_	1/14/202	1
Location	Product Name/Description	Size (oz)	Condition *	Chemical Ingred	ients	PID Reading	COC Y/N?
	See attached hand notes						
				h 1	***		
	,						
2							

			-	
Product Inventory Complete?	Yes	Were there any elevated PID readings taken on site?	Yes	▼ Products with COC?

^{*} Describe the condition of the product containers as Unopened (UO), Used (U), or Deteriorated (D)

^{**} Photographs of the **front and back** of product containers can replace the handwritten list of chemical ingredients. However, the photographs must be of good quality and ingredient labels must be legible.



Site Name: Bristol-Myers Squibb	Site Code:	Operable Unit:
Building Code: Industrial Building Name:	Building 3	
Address: 6000 Thompson Rd	Apt/	/Suite No:
City: East Syracuse State: N	Zip: 13057	County: Onondaga
Factors Affecting Indoor Air Quailty	NOTIFICATION OF THE STATE OF TH	Management of the Control of the Con
Frequency Basement/Lowest Level is Occupied?: ALMOST NEVER	Floor Material: CEM	MENT
☐ Inhabited? ☐ HVAC System On? ☐ Bath	room Exhaust Fan?	Kitchen Exhaust Fan?
Alternate Heat Source: NONE	☐ Is there	e smoking in the building?
Air Fresheners? Description/Location of Air Freshener:	Bathrooms	
Cleaning Products Used Recently?: Description of Cleaning Products:	Lysol and other	disinfectants
Cosmetic Products Used Recently?: Description of Cosmetic Products:	<u> </u>	
New Carpet or Furniture? Location of New Carpet/Furniture:	****	
Recent Dry Cleaning? Location of Recently Dry Cleaned Fabrics:	Various staff us	e dry cleaned clothing
Recent Painting/Staining? Location of New Painting:		
Solvent or Chemical Odors? Describe Odors (if any):		
Do Any Occupants Use Solvents At Work? If So, List Solvents Used:	Various for pipe f	ittings
Recent Pesticide/Rodenticide? Description of Last Use:		
Describe Any Household Activities (chemical use,/storage, unvented appliar The building is the boiler and maintenance area for has an exhaust fan, as well as a bathroom with clestock room have been identified, marked out, and I	nces, hobbies, etc.) That Ma r the facility. Th aners. The cleanin	ay Affect Indoor Air Quality: e water testing area g products in the
Any Prior Testing For Radon? If So, When?:		
Any Prior Testing For VOCs? If So, When?: Jan 1, 2019		
Sampling Conditions		
Weather Conditions: PARTLY CLOUDY Out	door Temperature:	°F
Current Building Use: MULTI-USE BUILDING Bard	ometric Pressure:	in(hg)
Product Inventory Complete? Yes Building Questionnaire	Completed?	



Site Name: BMS Syracuse Site Code: Operable Unit:
Building Code: Bulding Name: Building Name: Building Name: Building Name: Building Name:
Address: Apt/Suite No:
City: Sylvacure M State: MZip: County: USA
Factors Affecting Indoor Air Quailty
Frequency Basement/Lowest Level is Occupied?: 3 shifts/Ddy Floor Material: (oncete w/tile
Inhabited?
Alternate Heat Source: Flow Heater (Electric) Is there smoking in the building?
Air Fresheners? Description/Location of Air Freshener:
Cleaning Products Used Recently?: Description of Cleaning Products: Champer Spay Dighted A
Cosmetic Products Used Recently?: Description of Cosmetic Products:
New Carpet or Furniture? Location of New Carpet/Furniture:
Recent Dry Cleaning? Location of Recently Dry Cleaned Fabrics:
Recent Painting/Staining? Location of New Painting:
Solvent or Chemical Odors? Describe Odors (if any): Hand Son to
Do Any Occupants Use Solvents At Work? If So, List Solvents Used: Yes > See product in vertical
Recent Pesticide/Rodenticide? Description of Last Use:
Describe Any Household Activities (chemical use,/storage, unvented appliances, hobbies, etc.) That May Affect Indoor Air Quality: () Man' grand 4, FO: FACOAIO, Suple FD: Job 1 + 1A; Locatus; Bldg 3 Control Ross Vs AMIS ID: FACOAIS, Suple ID: Job 2 + 2M, Locatus; Bldg 3 Control Ross Vs Bldg 2 ID: FACOAIS, Sample ID: Job 3 + 3A, Locatus; Bldg 3 Control Ross Vs 3106 Drain
- Note Differential pressur Manutes Collected @ Cordnel Rom VS Blog 2, VS AMB, VS Dian.
Any Prior Testing For Radon? If So, When?:
Any Prior Testing For VOCs? If So, When?: 2019
Sampling Conditions Start Tenporate: 70.4 = IA Hamilty: 29.6
Current Building Use: Boiler Contai Room Outdoor Temperature: 40.8 7 62.8 % Hum °F Barometric Pressure: 5tart 29.8 4 in(hg)
Product Inventory Complete? 1/12/21 \(\mathbb{N}\) Building Questionnaire Completed? The End: Stop: Further top: 70.1 F 29.1% Hung



	PRODU	JCT INVENTORY		
Building Name: Building		Bldg Code:	Date: 1/12/2	1
Bldg Address:	BMS Syracuse		Apt/Suite No:	
Bldg City/State/Zip:	J			<u> </u>
Make and Model of PID:	.7 ev MiniKAE 300	U 10# FA0032	Date of Calibration: 1/12/21	@0930 KM

	Location	Product Name/Description	Size (oz)	Condition *	Chemical Ingredients / Phulo	PID (Pf Reading	COC Y/N
	B2B-C1	3-in-1: Mult; purpose Bil	3°Z	U	Petrolian distillates - 2104 -2105	0.0	
	11	Loctite-567 Threed Seglant	1.690Z *3	U	- 2106	0.0	
		Loctite - Gasket eliminator 518	10.15 02	U	- 2106	0.0	
	11	Loctite - 55 pipe Sealing Cord	~30Z	U	3:11can Fesia - 2106	0.0	
	11	Loctile - TRC/680 retaining compound	1,69 oz	U	-2108	0.0	
	11	Locite LB-8008	8°Z	U	Quartz - 2108 - 2109	0.0	
100	· ·	LA-CO: Slic-tite Thread Sealing compound	8ºZ	U	-2108	0.0	
	ų.	LA-CO: Epoxy Stick	~80Z	U	Bisphenol-A- Epoxy resin, -21 5:1:00 D'ox'de		
	11	Techspray-Transitur Siliene grease	402	U	- 2110		П
1	ĮĮ.	Blastic: Penetrating catalyst	11.z	U	Petrolium distillates -2110	4	П
1	62B-A1	Sodium Hydroxide 6	5 5 gallon>	U	-2113 -2114	1 ().()	П
1	82B-AZ		55 99/10ms	UD	- 2115 - 2116		П
1	32B-A2	Cooling Water treatment	~55 gallow	UO	Sodium hydroxide -2117 -2118		П
P	32B-A2	1 1 1 1 1 10	~55	UO	Sodium bromo sulfamate - 2119 Sodium Chloro sulfamate - 212		
B	2B-A2	Boiler water treatment BL122	120	U	Sodium Bisulfite -212	0.0	П
e	52B-A2	Boiler vater treatment BL 2452	120	U	Potasina Hydroxide -212 -212	3 0 0	

^{*} Describe the condition of the product containers as Unopened (UO), Used (U), or Deteriorated (D)

* Describe the condition of the front and back of product containers can replace the handwritten list of chemical ingredients. However, the photographs of the floor quality and ingredient labels must be legible.

priotograpiis			
Product Inventory Complete?	,	Were there any elevated PID readings taken on site?	Products with COC



CABIT	01 –				
	and a street of	PRO	DUCT INVENTORY	The second secon	W 2000 A 20
Building Name: Build	ing 2A -	Floor 1	Bldg Code: B2	A - C2 Date:	1/12/21
Bldg Address:	BMS	ATT OF THE PARTY OF	126	Apt/Suite No	0:
Bldg City/State/Zip:	The second second	J	A Design		
Make and Model of DID	1174	10.0	3000	Date of Calibration:	1/12/21

Location	Product Name/Description	Size (oz)	Condition *	Chemical Ingredients/Photo	PID Reading	COC Y/N?
32A-C2 Top Shif	CRC-HF Contact	110Z	U	Petrolium distillates - 2125 - 2126	0.0	
II A	CRC - 2-26	Noz	U	Potoliam distillates -2127 -2128 -2129	0.0	
11	Sprayon - Multipurpuse adhesive	16.75 0Z	U	Hexane, Actone, propare, -2125 Bytane	0.0	
ıl 🥳	Flex Seal Liguid Rubber Coating	1402	U	Toluene, Minural spirits, -2125 petrolium Solvent -2131	0.0	
lr_	Flex Seal light	1280Z ×2	U	-2132 -2135 -2136	3.3	
11	Sprayon 300601 Red insulating Varnish	15.25 oz	U×I	Xylene, Actore, tolure, 1-2134 propore, Butane, iron pxide	0.0	
11	Krylm - Rost tough	150Z	UXIO	-2134	0.0	
11	Cal-SPRAY Coil cleaner	18.z	U ->	Trichloroethylere - 2134 -2137	0.0	¥
11	NAPA-10W-30 Motor Oil	32oz	U	-2138	0.0	П
11	Christerton - 723 Spranolus - pertutino	12.30Z	U	-2138	0.0	П
1)	E-Z Weld 914 Dipe Cleaner	802	U	Authyl Effyl Ketone -2138 -2139	0.0	Г
11	"E-Zweld" PVC count	1602	U×2	Tetahydroturan -2138 MEK -2140	1.9	Г
11	Carlon - Clear Primer	3202	7	MEK, Cyclohexanore, -2141 Tetrahydrofuran, -2142	429	
11	Spray On - Molly Chain Lube	1102	U×2	Propare, herare, -2141 -2143	1.5	П
11	Spruzon - Anti Sieze Lube	11.25 02	U	2141	0.0	П
- 11	Rustoleum - Enamel	15.Z	U×6	Actor, Ethyl Benzere, -2141 Xyline -2145	0.8	

Product Inventory Complete?	Were there any elevated PID readings taken on site?	Products with COC
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^{*} Describe the condition of the product containers as **Unopened (UO), Used (U)**, or **Deteriorated (D)**** Photographs of the **front and back** of product containers can replace the handwritten list of chemical ingredients. However, the photographs must be of good quality and ingredient labels must be legible.



Structure Sampling Questionnaire and Building Inventory

New York State Department of Environmental Conservation

Building Name	e: Building 2A -		ODUCT INV	Code: B2 A - [C2] Date:	1/12/	121
Bldg Address:	- Daniona En		racuse	Apt/Suite N	I MADE TO THE	
Bldg City/Stat		J	OF ELECTRICAL MARKET			
Make and Mo	del of PID: 11, 7 eV	Mini K	lae 301	Date of Calibration:	1/12/	21
Location	Product Name/Description	Size (oz)	Condition *	Chemical Ingredients / Photo	PID Reading	COC Y/N
B2A - C2 Middle shelf	Virginia - Vaccume	1 Quart	U	- 2146	0.0	
N _	Chemsafe - Ram truck-Sprun Adksive	120Z	U	n-butane, propone, -2146 Acetone -2147	0.4	
u	CROWN - Spray	14.102	U	Actor - 2147 Actor, heptone, -2148 propone -2148	0.5	
11	Sprayon - Chain + Pin bushing lube	1102	D	-2146	0.0	
11	Hercules - OCIERT PV C PRIMER	16 02	U	Tetrahydrofuran, (yeluhour, 1-215) MEK, Acetone -2151	0.0	
h	Steel IT - Anti Rust Stainless Steel	1402	Vo	Polyur ethane, xyol, Dress, Aliohatic Hadrocardon -2152	0.0	
h	Chisterton 730 SPRAGRIP	12 oz	U	Isobutan, -2150 huptone -2153	0.0	
11	Crown - TFE	10,5	U	Xylene, Active, 1-215 Cycloberac, Dimethyl ether 1-215	0.5	
11	CAL - SPRAY instant Irah Sealer	1602	U	-2150	0.7	
11	Virginia auto #10	1 gal	U	Aceton, tetrachors et hylen, 1-215 Hydro Carbons 1-215	13.9	
11	3M Nitrile pastice	5°z	UXI UUX2	- 2156 - 2158	0.0	
B2A-C2 Button	SW-DTM ACOULIC	1991	Ux3	-2160 -2161	0.0	
11	SW- usethane Alkyd	Igal	U Alter	- 2160 - 2162	0.0	
11	Temper Hote - high	lyal	U	Tolyane Petrolium Distillates -2163	0.7	
-11	SW-A100 8 Acrylic Formula	5gal	U	-2164 -2165	0.0	
11	Sunny Side-Low odur mineral spirits	Igal	U×2	Petrolium -2164 distillates -2166	0.2	П

Product Inventory Complete?	Were there any elevated PID readings taken on site?	Products with COC
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^{**} Photographs of the front and back of product containers can replace the handwritten list of chemical ingredients. However, the photographs must be of good quality and ingredient labels must be legible.



	State/Zip: Model of PID:!!, 7e V	Mini Ra	e 300	Date of Calibration:	1/2/2	1
Locatio	n Product Name/Description	Size (oz)	Condition *	Chemical Ingredients / hch	PID Reading	cc
B2A-C Bottom	2 Enerpac Hydraulic	1gal	U	- 2164	O.O	
BZA-C Top shift	3 Snoup-leak detecto	-	041 041	-2169	0.0	
'n	Weld-on-PC64 Primer Cleaner	32°Z	D	Actore, MEK, 1-2169, -2170 tetrahydro Furan, Cyclubusan	0.6	
11	Rushleum Primer	Igal	υ	-2169 -2171	0.0	
11	11rglon - Iron gurd Accylic	الإما	U×3	-2169 -2172	0.0	
ч	PPG - Multiprime Alkyd primer	Igal	U	-2174 - 2175	0.0	
11	Hercules - PVC Primer - Clear	32°z	U×2	Aceton, CycloHexan, MEK, -217 tetrahydrofuran -2172	4 00	
11	Hercules - PUC+ CPUC Primer - purple	3202	U	1 -2174	-	
μ	Hercules - PUC - plastic pipe cement	3202	U	-2174 -2179		
1	Herchles - PVC -	at 32.2	U	· -2174 -2179	12	
įı	EZ-Weld Parple Pr	imer 1602	U	Tetra Alydro Flura, MEN, 1-218 Cyclobrane, Acetone -218	0 00	
11	150 - Heet injector cleans	1202	U	150 propy 1 Alcoh.) -2181 -2181	0 00	
1)	Virginia - Vaccine Dump oil	lquet	U	-218		
11	NAPA - Motor oil 5W-30	1 gourt	U × 1	- 2180	0.0	
BZA-C3 middle st	EMMarate - Refrigerat	m I gal	U×2	Polyol ester - 2183 - 2184	3 00	
11	Com Star - Co. 1 flus Cleaner - CF20	h 1991	U	- 21 83		



	Make and Model of PID: 11.7 eV Min: Rae 3000 Date of Calibration: 1/12/21									
	Location	Product Name/Description	Size (oz)	Condition *	Chemical Ingredients	PID Reading	COC Y/N			
B2A- C3	B2A-C3 middle shelf		1971	Ux2	Mineral 0:1 -2183 -2185	0.0				
	11	Virginia AB150	Igal	U	-7183	0.0				
2	11	SW-12dustrial	Igal	Ux3	6+hyl Denzene -2186 -2187	0.0				
	11	Virginia 10n Degressing Sulvent	Igal	UXI VOX3	Accton, tetral Moroethylen, 1-2186 Hydrocarbon 1-2188	0.2				
	11	SW- Promer 200	1941	U×I	-7186 -2189	0.0	Ō			
62A-C3	B2A-C3 Buttom	Edmac - Synthetic	5gal	U	- 2190	0.0				
· C3	11	TRANE - 0:122 Refridgeration oil	2.5 59/	U	- 2190	0.0				
	B2A-AI	DowFrost	55gal	U×2	- 2191 - 2192	0.0				
A	lı	Permutrent PC191T	5gal	U×2	-2194	0.0	П			
1	11	Nalsperse 73550	5gal	U×2	-2194	0.0	П			
	li	Nalco 7469 Antiform	5941	U	-2194	0.0				
	11	Nalco 7341	5991	U	- 2195	0.0				



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PRODUCT INVENTORY

Building Name: Building 2A - Floor | Bidg Code: B2A - [A2+A3] Date: 1/14/21

Bldg Address: BMS Syracusc Apt/Suite No:

Bldg City/State/Zip:

Make and Model of PID: 11.7 eV Min'i Race 3000 Date of Calibration: 1/14/21

Location	Product Name/Description	Size (oz)	Condition *	Chemical Ingredients	PID Reading	COC Y/N?
B2A-A2	Chem treat - Coul water treatment CL683	55ga)	UXG	Sodium hydroxide - 2198	0.0	П
II.	Chen treat-CL49	55 gal	U×6	-2199 -2200	0.0	
į)	Chun treat CL2150	55	Ux6	-2201 -2202 -2196	0.0	
н	Ant: scalant	85	U	-2196	0.0	
1	Sodium Bisulfite	75	U	-2197	0.0	
B2A-A3	Chantrent- Cul water treatment CLG 836	55gn/	U	Sodium hydraide	0.0	
11	Chentrat CL49	55gal	U	-7245	0.0	
4	Chentrent CL2150	55gal	U	12.145	0.0	
	Page Total Arriva	5-1		-7208	0.0	П
	1 to the of phase or	124	10.05	-72.04	33	П
	RED TOW 30	12 /		- 7. % 100	8 0	
	mad 50.224	8),	9 () ((4))	2 Lon	94	П
	The state of the state of	190	-0	27.07	0.0	П
	and Hard State	1 6 m	V	Marylamorm - 2207	0.7	
						П

* Describe the condition of the product containers as Unopened (UO), Used (U), or Deteriorated (D)

** Photographs of the front and back of product containers can replace the handwritten list of chemical ingredients. However, the photographs must be of good quality and ingredient labels must be legible.

Product Inventory Complete?	Were there any elevated PID readings taken on site?	Products with COC
Ploduct machinery company		Floudets with Coc



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Structure Sampling Questionnaire and Building Inventory New York State Department of Environmental Conservation

		PR	ODUCT INVENTO	DRY
Building Name: Buildi	nd 3 -	Floor	Bldg Code:	B3 - [Lab+BR] Date: 1/14/21
Bldg Address:		Syracun		Apt/Suite No:
01d- Cla-15		J		A STATE OF THE STA
Bldg City/State/Zip:				

	Location	Product Name/Description	Size (oz)	Condition *	Chemical Ingredients / Photo	PID Reading	COC Y/N?
	B3-Lab	EMD-Buffer Solution OH 4.0	500 mL	U	-2203	0,0	
	u	EMD-Buffer Solution	500 ML	U x 2	- 2203	0.0	
	10	EMD-Buffer Schoten	500	U *2	- 2203	0.0	
	11	Nalco-ORP Stenderd	1 4	U×2	-2204	0.0	
53	11	Nalco - ORP stander	11	UO	- 2204	0.0	
N-LA	11	Nalco - 50980 - Trasar 3 Cal	1L	U	-2204	0.0	
B	11	Nalco - 50920	11	U×1	-2205	0.0	
	11	Nalco - SOLN ST-2 doponixe sulfite	11	U×2	- 2205	0.0	П
	11	Nalco Total Alkadiata		U×2	-2205	0,0	
	11	NAICO - Phenol phylein	500mL	U×1	-2206	0.3	П
)(NALCO - Traser 30	4 L	UXI	-2206	0.0	
	11	NALCO - SOLN 50226 H2SO4 Titrant	4L	U × I	- 2206	0.0	0
F7	83-BR	Tough Guy- hand cleaner	1991	U	- 2207	0.0	П
- BR	11	Champion - Sprandibafed as	0.50z	U	Dinethyl ammonium -2207 chloride	0.0	
						1	

^{*} Describe the condition of the product containers as Unopened (UO), Used (U), or Deteriorated (D)

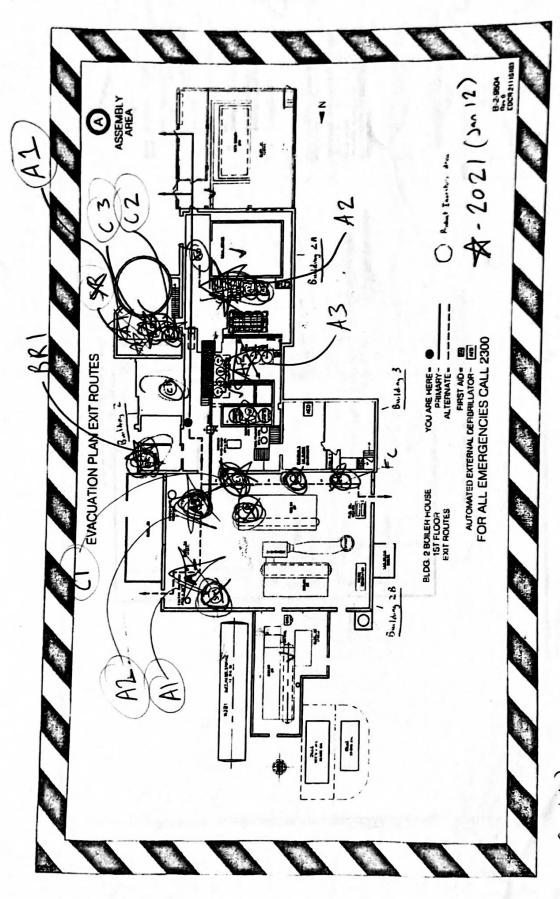
Product Inventory Complete?	Were there any elevated PID readings taken on site?	Products with COC

^{**} Photographs of the **front and back** of product containers can replace the handwritten list of chemical ingredients. However, the photographs must be of good quality and ingredient labels must be legible.

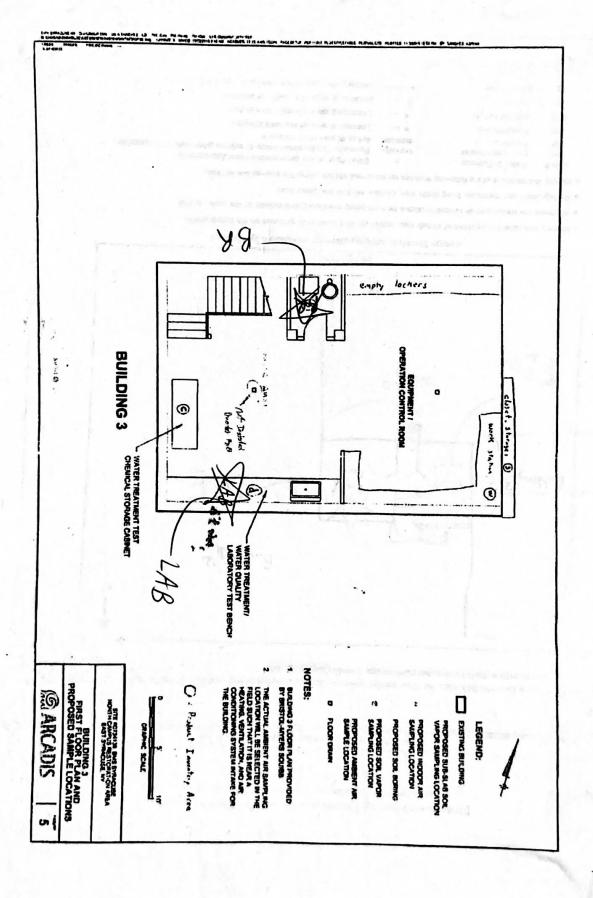


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	A 13 A =	PI	RODUCT INVEN	TORY		101
Building Name	E. Building 3 - Flu	ur 2	Bldg Code	B3-F2[E . W. bote:		21
Bldg Address:	Building 3 - Flu	Syra	cuse	Apt/Suite N	Ot	
Bldg City/Stat	e/Zip:	U				
Make and Mo	del of PID: 11.7 cv 1	Mini Rac 3000 Date of Calibration: 1/14/				
Location	Product Name/Description	Size (oz)	Condition *	Chemical Ingredients	PID Reading	COC Y/I
B3-F2E	Virginia - Alk: Foam	Igal	U×2	- 2208	0.0	
n	Bright DYES	194	U×2	-2209	0.0	п
tı	TRANE - 01200022	2.5	UxI	- 2210	0.0	П
	0					П
						П
						П
						0
						П
						0
						0
	*					D
					1	D
			177			1 0



28-C1 24.C2





FIRST FLOOR BUILDING LAYOUT SKETCH Please click the box with the blue border below to upload a sketch of the first floor of the building. The sketch should be in a standard image format (.jpg, .png, .tiff) Clear Image Building 2 Building 3 Bulling 3 Design Sketch Design Sketch Guidelines and Recommended Symbology a identify and tabel the locations of all sub-stab, indoor air, and outdoor air samples on the layout sketch. a Measure the distance of all sample locations from identifiable features, and include on the layout sketch. a identify room use (bedroom, living room, den, kitchen, etc.) on the layout sketa identify the locations of the following features on the layout sketch, using the appropriate symbols BOF Boiler or Furnace Other floor or wall penetrations (label appropriately) Penmeter Drains (draw inside or outside outer walls as appropriate HW Hot Water Hester Fireplaces ****** Areas of broken-up concrete WS Wood Stores Location & label of sub-sizb samples 0 SS-1 WID Washer / Dryer Location & label of Indoor air samples Sumps 0 Cat. 1 Location & tabel of outdoor air samples Floor Drains • FFET-1 Location and label of any pressure field test holes Product Investig Ann



Building Code:	m/Indestrial	Address: Bailding	13		
Sampling Inform	ation				
Sampler Name(s):	Dayiel Zuc	K	Sampler Con	npany Code: <u></u> <i>Aira</i>	ndis
Sample Collection Da				es Sent To Lab:	
Sample Chain of Cust	tody Number:	_		Sample Location ID:	4
SUMMA Canister	Information				
Sample ID:	IA-3 (DUP-0/1	1421) DUP-011421	AMB-011421		
Location Code:	Indoor Air	IA	AMB-		
Location Type:	(ummella)	(annocial	Conherical		
Canister ID:	6L0764	2290	640505		
Regulator ID:	23485	24335	23522		
Matrix:	A-Air	JA:AIr	Ambiait Air		
Sampling Method:	6L Symmo - 70-15	61541ma -70-15			
Sampling Area In	fo				
Slab Thickness (inches):	M	M	M		
Sub-Slab Material:	M	MA	M		
Sub-Slab Moisture:	M	M	MA		
Seal Type:	M	MA	M		
Seal Adequate?:	$\mathcal{N} \Box$	NA [MAD		
Sample Times and	Vacuum Readings				
Sample Start Date/Time:	0840 1/14/21	0840 1/14/21	0838 1/14/21		\
Vacuum Gauge Start:	-27.5 ([re: -30.20)	-27.5 (30.16)	-28.75 (Pre-2059)		
Sample End Date/Time:	1642 1/14/21	1642 1/14/20	1640 1/4/21		
Vacuum Gauge End:	-6.25	-6.75	-6.5		
Sample Duration (hrs):	Shy	8 his	8 hs		
Vacuum Gauge Unit:	inHg	in Hg	in Hy		
Sample QA/QC Rea	adings				
Vapor Port Purge:	M	M	M		
mini RAE 3000 11.7 Lap	A: 0.0 ppn	0.0 ppm	0.0 ppm		
Purge PID Unit: 3	U. U spin	O. Oppm	0.0 ppm		
Tracer Test Pass:					. 🗆
Sample start	and end times should	be entered using	the following form	at: MM/DD/YYYY	HH:MM



Indoor/Ambient Air Sample Collection Log

		Sample ID: 14-3		
Client:	BMS	Date/Day:	5/23/2019 4/29/21	
Project:	Buildy 3 VI	Sample Intake Height:	4' ALS	
Location:	Boiler Control Room	Subcontractor:	NA	
Project #:		Miscellaneous	-1-11	
Samplers:	Daniel Zuck	Equipment:	IA Tracknet and Nor Air 800	
Coordinates:	(See attached Figure)	Time Start:	0805	
Outdoor/Indoor:	Ludoov	Stop Tine	1610	

Instrument Readings:

Time	Canister Pressure (inches Hg)	Temperature (F)	Relative Humidity (%)	Air Speed (ft/min)	Barometric Pressure	PID (ppb) / (ppm)
0805	-29.5	69.4	49.7	0	29.37	0
1345	-10	73.3	48.6	0	29.23	
1610	-6	74.2	51.5	0	29.17	0

SUMMA Canister Information

a.	/ · ·		
Size	circ	e one):	

Canister ID:

6L0742

Flow Controller ID: 23407

Photos: 2021.04.29 _ 081036	
=DA Sucien Plus Pole P. Efflort 25t ft/	
= DA System flow Role @ Effloot 255 ft/mm = Staff were using hand sometizing sping	
50 7	



Indoor/Ambient Air Sample Collection Log

		Sample II	D: T#-3
Client:	BM5	Date/Day:	5/23/2019 (9) 5/6/21
Project:	13/15 Building 3	Sample Intake Height:	4' ALS
Location:	Busement antici ilcom	Subcontractor:	NA
Project #:		Miscellaneous	TA Calpa track
Samplers:	Daniel Zuck	Equipment:	TA Caba tracket
Coordinates:	(See attached Figure)	Time Start:	0825
Outdoor/Indoor:	Inlege	time Ston	1632

Instrument Readings:

Time	Canister Pressure (inches Hg)	Temperature (F)	Relative Humidity (%)	Air Speed (ft/min)	Barometric Pressure	PID (ppm)
0825	-30	70.0	45.2	0	79.64	0
0941	-24	70,2	41.9	0	29.65	6
1632	-6	71.2	33.5	0	29.60	0

SUMMA Canister Information

Size (circle one):

1 L 6 L

Canister ID:

Flow Controller ID: 25317

Photos: 2021 0506 083511.jpg	
Average (FM - 279	



Indoor/Ambient Air Sample Collection Log

		Sample ID:	IA- 3
Client:	BMS	Date/Day:	5/13/21 74.
Project:	Building 3 VI	Sample Intake Height:	4' ALS
Location:	Syvacuse, NY	Subcontractor:	NA
Project #:	, ,	Miscellaneous	
Samplers:	D. Zuck	Equipment:	Novan 800
Coordinates:	(See attached Figure)	Time Start:	0815
Outdoor/Indoor:	IA	End Time:	1615

Instrument Readings:

Time	Canister Pressure (inches Hg)	Temperature (F)	Relative Humidity (%)	Air Speed (ft/min)	Barometric Pressure	PID (ppb)
08/5	-31	\$ 69.9	35.1	0	29.87	0
1615	-6	71.6	31,7	0	29.80	Ò
	,					

SUMMA Canister Information

Size	circ	le	one)	:

Canister ID:

6L0453

Flow Controller ID: 23255

Photos: 2021 3513, 1625 38	
Average 251 CFM our limin data Colletion time	
	123101



Indoor/Ambient Air Sample Collection Log

		Sample ID:	IA-3
Client:	BM5	Date/Day:	5/20/21 thursday
Project:	Buddy 3 FA	Sample Intake Height:	4' ALS
Location:	Boiler Control Room	Subcontractor:	NA
Project #:		Miscellaneous	
Samplers:	D. Zuck	Equipment:	Novid 800
Coordinates:	(See attached Figure)	Time Start:	0828
Outdoor/Indoor:	IA	End Time:	1628

Instrument Readings:

Time	Canister Pressure (inches Hg)	Temperature (F)	Relative Humidity (%)	Air Speed (ft/min)	Barometric Pressure	PID (ppb)
0828	-50.52	70.0	50.9	0	29.99	0
1628	-6	73,45	54,2	0	29.92	0

SUMMA Canister Information

Size (circle one):

1 L 6 L

Canister ID:

6L2210

Flow Controller ID:

23477

Photos: 20210520_215346			
	1545 3/550		
Eftbert -1 Sayle Collected a	0345 Calb: 660902/F	C#: 23828* PV	19560: -21.5/-5
Int/104-1 1	161071615 114; GL07359 F	(H: 73466*	· · · -28.5 /->
* FL's had flow Regulators Renoe	1 to collect quick grob samples	from port.	
		100	

ARCADIS

Indoor/Ambient Air Sample Collection Log

	A CONTRACTOR OF THE PARTY OF TH	Sample ID:	AA- 3
Client:	BMS	Date/Day:	Thorday C/24/21
Project:	Bailding 7 VI	Sample Intake Height:	4' ALS
Location:	Stracuse MY	Subcontractor:	NA
Project #:	300 649 43.00001	THE NAME OF STREET AND ADDRESS.	r Street - No.
Samplers:	D. Zuck	Equipment:	Vehicle Porking
Coordinates:	(See attached Figure)	Time Start:	1030
Outdoor/Indoor:	A A		A STATE OF THE PARTY OF THE PAR
	Au .	End Time:	1830

Instrument Readings:

Time	Canister Pressure (inches Hg)	Temperature (F)	Relative Humidity (%)	Air Speed (ft/min)	Barometric Pressure	PID (ppb)
1030	-30	76.64	35,22	0	30.34	0
1430	-15.5	88.10	25.11	0	30.29	0
1430	-7	85.60	28.40	0	30.23	0
	yr. 87 Kiper, m					
700						
			# 4 Alex			
				and the second		

SUMMA Canister Information

Size (circle one):

1L 6L

Canister ID:

661321 23401

Flow Controller ID:

Photos:	Ihr	pressure	AA - 26.5	 and the second s	
@	446	pressing	AA - 15.5	the state of the s	100
					And the second

ARCADIS

Indoor/Ambient Air Sample Collection Log

	American Control of Co	Sample ID:	IA-3 / DUP-06242
Client:	BM5	Date/Day:	Thusday 6/24/21
Project:	Building 3VI	Sample Intake Height:	4' ALS
Location:	Syracuse NY	Subcontractor:	NA
Project #:	Syraluse NY 30064943.00001	Miscellaneous	The second secon
Samplers:	D. Zuck	Equipment:	Norther 800
Coordinates:	(See attached Figure)	Time Start:	1025
Outdoor/Indoor:	IA	End Time:	18 25

Instrument Readings:

Time	Canister Pressure (inches Hg)		Temperature (F)	Relative Humidity (%)	Air Speed (ft/min)	Barometric Pressure	PID (ppb)
W25	- 30	-30	69.1	40.71	0	30.34	0
1430	-15.5	-15	74.2	46.85	0	30.29	0
1825	-6,0	-6.0	77.09	43.31	0	30.24	0
			- facility	A		Total Control of the	
		9	1	A war		+	- 1 + 1
			1	NAME BY AND ADDRESS OF THE PARTY OF THE PART			

SUMMA Canister Information

Size	Simo	A AM	•1•
		ic vii	

1L 6L

GLD384] Dul

Canister ID:

66038 6L0062

Flow Controller ID: 23593 23319

Photos:	
- @ 1 hr Presson: If = 265/	Dal-57
- Pyhr Passue : IA/.	15.5 / Dup -15.0



Building 3 Vapor Intrusion and Air Treatment System Assessment Report Bristol-Myers Squibb Syracuse North Campus Restoration Area (Site #C734138) East Syracuse, New York



Building 3 Air Treatment System and Pilot Testing



Building 3 Vapor Intrusion and Air Treatment System Assessment Report Bristol-Myers Squibb Syracuse North Campus Restoration Area (Site #C734138) East Syracuse, New York



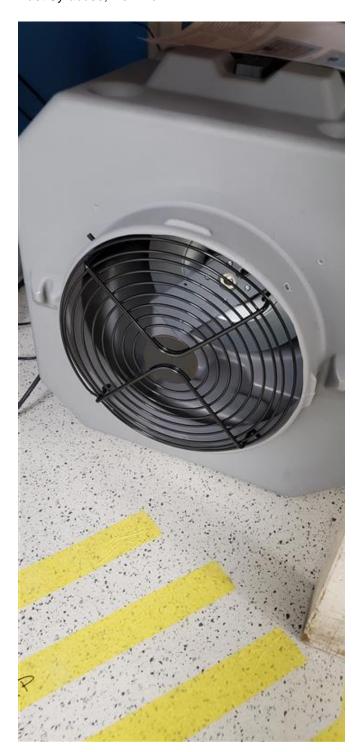


Building 3 Air Treatment System

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Building 3 Vapor Intrusion and Air Treatment System Assessment Report Bristol-Myers Squibb Syracuse North Campus Restoration Area (Site #C734138) East Syracuse, New York





Building 3 Air Treatment System

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Building 3 Vapor Intrusion and Air Treatment System Assessment Report Bristol-Myers Squibb Syracuse North Campus Restoration Area (Site #C734138) East Syracuse, New York

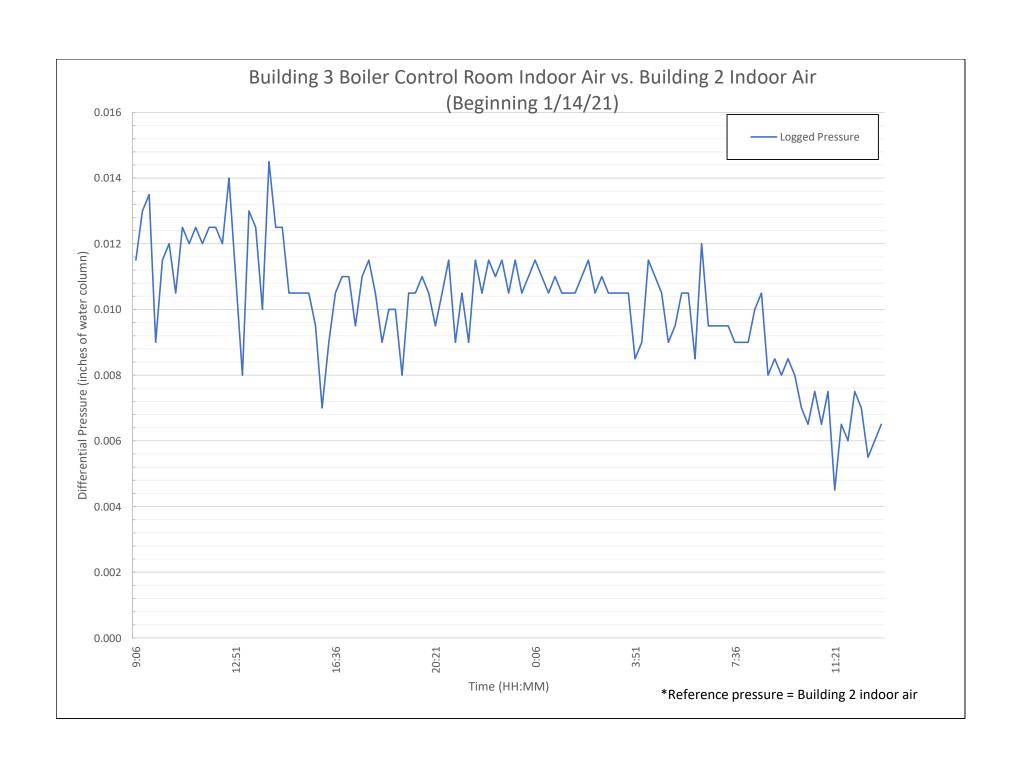
Air Sampling Locations

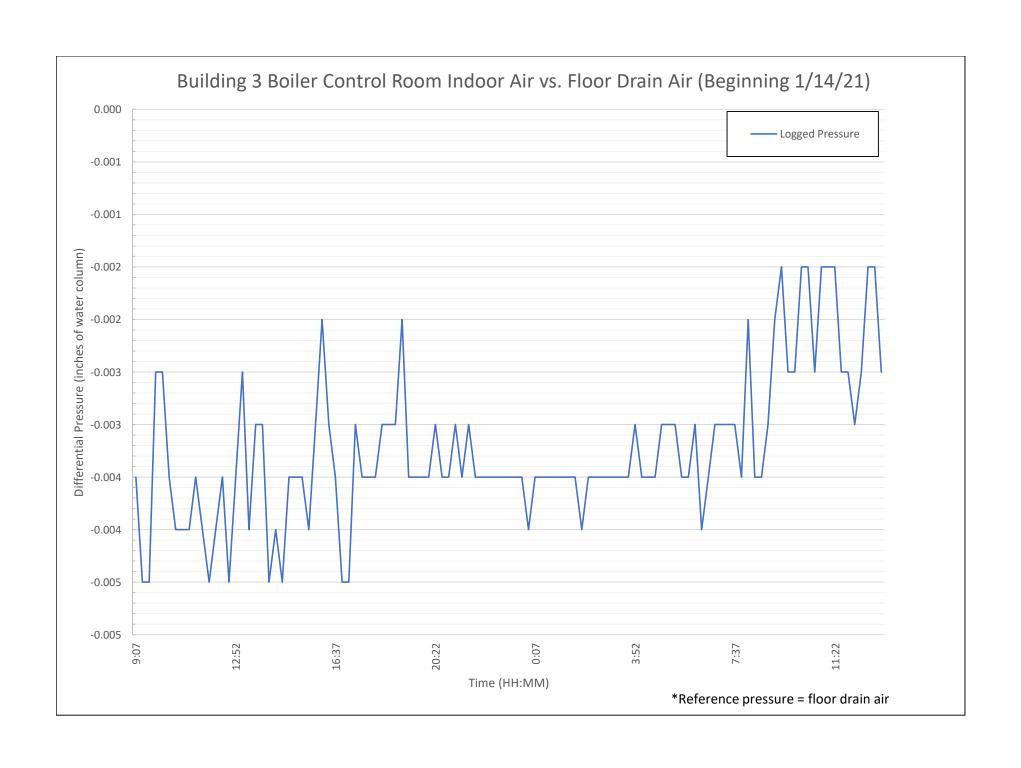


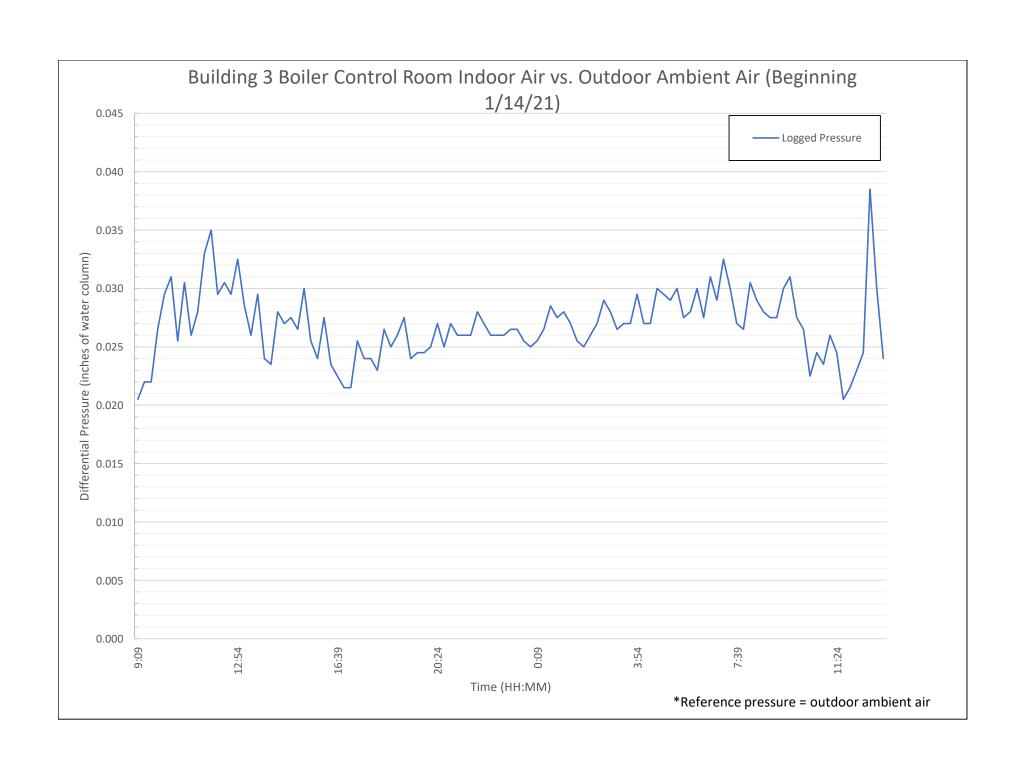


Appendix B

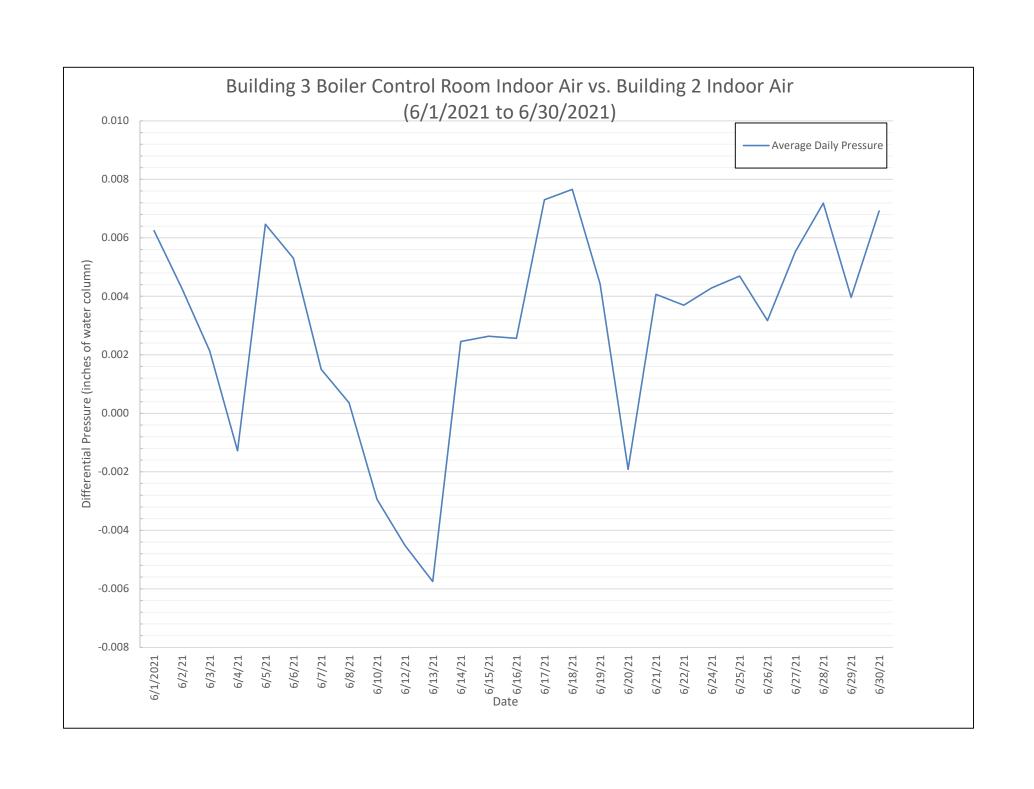
Differential Pressure Graphs

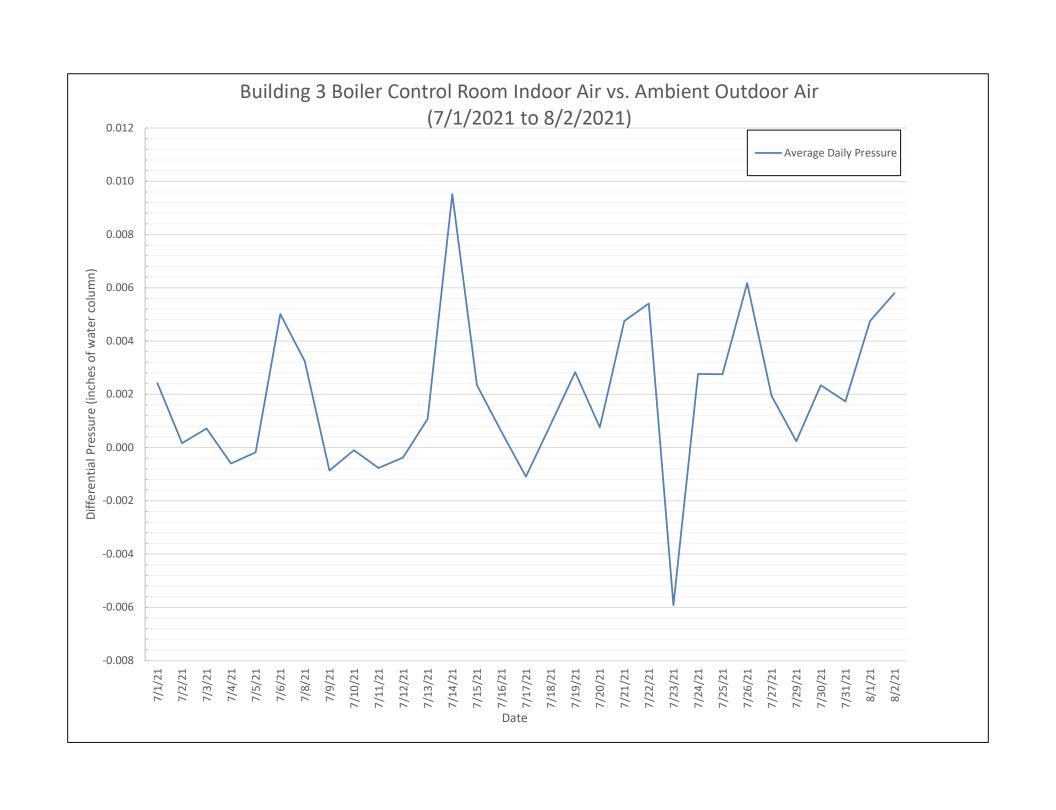












Appendix C

Data Usability Summary Reports and Laboratory Report Data



Bristol Myers Squibb Thompson Road Investigation

Data Usability Summary Report

Syracuse, NY

Volatile Organic Compound (VOC) Analysis

SDGs # 2105446

Analyses Performed By: Eurofins Air Toxics

Folsom, CA

Report #41878R Review Level: Tier III Project: 30064943

SUMMARY

This data quality assessment summarizes the review of Sample Delivery Group (SDG) # 2105446 for samples collected in association with the Bristol Myers Squibb Thompson Road Site. The review was conducted as a Tier III evaluation and included review of data package completeness. Only analytical data associated with constituents of concern were reviewed for this validation. Field documentation was not included in this review. Included with this assessment are the validation annotated sample result sheets, and chain of custody. Analyses were performed on the following samples:

0 1 15			Sample	Parent		Į.	Analysis		
Sample ID	Lab ID	Matrix	Collection Date	Sample	VOC	svoc	PFAS	MET	ALD
IA-3	2105446-01A	Air	5/20/2021		Х				
Effluent-1	2105446-02A	Air	5/20/2021		Х				
Influent-1	2105446-03A	Air	5/20/2021		Х				

ANALYTICAL DATA PACKAGE DOCUMENTATION

The table below is the evaluation of the data package completeness.

	Performal Reported Acceptal			Not	
Items Reviewed	No	Yes	No	Yes	Required
Sample receipt condition		Х		Х	
2. Requested analyses and sample results		Х		Х	
Master tracking list		Х		Х	
4. Methods of analysis		Х		Х	
5. Reporting limits		Х		Х	
6. Sample collection date		Х		Х	
7. Laboratory sample received date		Х		Х	
8. Sample preservation verification (as applicable)		Х		Х	
9. Sample preparation/extraction/analysis dates		Х		Х	
10. Fully executed Chain-of-Custody (COC) form		Х		Х	
11. Narrative summary of QA or sample problems provided		Х		Х	
12. Data Package Completeness and Compliance		Х		Х	

Note:

QA - Quality Assurance

List of Acronyms

%D: Percent Difference

%R: Percent Recovery

AC: Acceptable

ALC/GLY: Alcohols/Glycols

BAL: Blank Action Level

CCV: Continuing Calibration Verification

CRDL: Contract Required Detection Limit

D: Dilution

EIS: Extractable Internal Standard

FB: Field Blank

FD: Field Duplicate

ALD: Aldehydes

GC/ECD: Gas Chromatograph/Electron Capture Detector

GC/MS: Gas Chromatograph/Mass spectrometer

HT: Holding Time

ICP: Inductively Coupled Plasma

ICS: Interference Control Sample

ICV: Initial Calibration Verification

ISTD: Internal Standards

LabDup: Laboratory Duplicate

LCS: Lab Control Sample

LCSD: Lab Control Sample Duplicate

LL: Lower Control Limit

MB: Method Blank

MDL: Method Detection Limit

MET: Metals

MS: Matrix Spike

MSD: Matrix Spike Duplicate

N/A: Not Applicable

NC: Not Compliant

List of Acronyms, Continued

PAH: Polyaromatic Hydrocarbon

PCB: Polychlorinated Biphenyl

PEST: Pesticide

PFAS: Per- and Polyfluoroalkyl Substances

QA: Quality Assurance

QC: Quality Control

RB: Rinse Blank

RL: Reporting Limit

RPD: Relative Percent Difference

RRF: Relative Response Factor

RSD: Relative Standard Deviation

RT: Retention Time

SDG: Sample Delivery Group

SerDil: Serial Dilution

SIM: Single Ion Monitoring

SOP: Standard Operating Procedure

SSTD: Surrogate Standards

SVOC: Semivolatile Organic Compound

TB: Trip Blank

TIC: Tentatively Identified Compound

TOC: Total Organic Carbon

TOTDIS: Total and Dissolved

UL: Upper Control Limit

USEPA: United States Environmental Protection Agency

VOC: Volatile Organic Compound

ORGANIC ANALYSIS INTRODUCTION

Analyses were performed according to United States Environmental Protection Agency (USEPA) Method TO-15. Data were reviewed in accordance with USEPA National Functional Guidelines of October 1999, USEPA Region II SOP HW-31- Validating Air Samples Volatile Organic Analysis of Ambient Air In Canister by Method TO-15 of October 2006, New York State DEC Analytical Method ASP 2005 TO-15 (QA/QC Criteria R9 TO-15), and NYSDEC Modifications to R9 TO-15 QA/QC Criteria October 2009.

The data review process is an evaluation of data on a technical basis rather than a determination of contract compliance. As such, the standards against which the data are being weighed may differ from those specified in the analytical method. It is assumed that the data package represents the best efforts of the laboratory and had already been subjected to adequate and sufficient quality review prior to submission.

During the review process, laboratory qualified and unqualified data are verified against the supporting documentation. Based on this evaluation, qualifier codes may be added, deleted, or modified by the data reviewer. Results are qualified with the following codes in accordance with USEPA National Functional Guidelines:

- Concentration (C) Qualifiers
 - U The compound was analyzed for but not detected. The associated value is the compound quantitation limit.
 - B The compound has been found in the sample as well as its associated blank, its presence in the sample may be suspect.
- Quantitation (Q) Qualifiers
 - E The compound was quantitated above the calibration range.
 - D Concentration is based on a diluted sample analysis.
- Validation Qualifiers
 - J The analyte was positively identified; however, the associated numerical value is an estimated concentration only.
 - UJ The analyte was not detected above the reported sample detection limit. However, the reported limit is approximate and may or may not represent the actual limit of detection.
 - UB Analyte considered non-detect at the listed value due to associated blank contamination.
 - N The analysis indicates the presence of a compound for which there is presumptive evidence to make a tentative identification.
 - R The sample results are rejected.

Two facts should be noted by all data users. First, the "R" flag means that the associated value is unusable. In other words, due to significant quality control (QC) problems, the analysis is invalid and provides no information as to whether the compound is present or not. "R" values should not appear on data tables because they cannot be relied upon, even as a last resort. The second fact to keep in mind is that no compound concentration, even if it has passed all QC tests, is guaranteed to be accurate. Strict QC serves to increase confidence in data but any value potentially contains error.

VOLATILE ORGANIC COMPOUND (VOC) ANALYSES

1. Holding Times

The specified holding times for the following methods are presented in the following table.

Method	Matrix	Holding Time	Preservation	Return Canister Pressure
USEPA TO-15	Air	30 days from collection to analysis	Ambient Temperature	< -1" Hg

All samples were analyzed within the specified holding time and canister return pressure / vacuum criteria.

2. Blank Contamination

Quality assurance (QA) blanks (i.e., method and rinse blanks) are prepared to identify any contamination which may have been introduced into the samples during sample preparation or field activity. Method blanks measure laboratory contamination. Rinse blanks measure contamination of samples during field operations.

A blank action level (BAL) of five times the concentration of a detected compound in an associated blank (common laboratory contaminant compounds are calculated at ten times) is calculated for QA blanks containing concentrations greater than the method detection limit (MDL). The BAL is compared to the associated sample results to determine the appropriate qualification of the sample results, if needed.

All compounds associated with the QA blanks exhibited a concentration less than the MDL, with the exception of the compounds listed in the following table. Sample results associated with QA blank contamination that were greater than the BAL resulted in the removal of the laboratory qualifier (B) of data. Sample results less than the BAL associated with the following sample locations were qualified as listed in the following table.

Sample Locations	Analytes	Sample Result	Qualification	
	Benzene			
Effluent-1	Ethylbenzene	Detected sample results <rl <bal<="" and="" td=""><td colspan="2" rowspan="2">"UB" at the RL</td></rl>	"UB" at the RL	
	m,p-Xylenes			

Note:

RL Reporting limit

3. Mass Spectrometer Tuning

Mass spectrometer performance was acceptable and all analyses were performed within a 12-hour tune clock.

System performance and column resolution were acceptable.

4. Calibration

Satisfactory instrument calibration is established to ensure that the instrument is capable of producing acceptable quantitative data. An initial calibration demonstrates that the instrument is capable of acceptable performance at the beginning of an experimental sequence. The continuing calibration verifies that the instrument daily performance is satisfactory.

4.1 Initial Calibration

The method specifies percent relative standard deviation (%RSD) and relative response factor (RRF) limits for select compounds only. A technical review of the data applies limits to all compounds with no exceptions.

All target compounds associated with the initial calibration standards must exhibit a %RSD less than the control limit (30%) or a correlation coefficient greater than 0.99 and an RRF value greater than control limit (0.05).

4.2 Continuing Calibration

All target compounds associated with the continuing calibration standard must exhibit a percent difference (%D) less than the control limit (30%) and RRF value greater than control limit (0.05).

Compounds associated with the calibrations were within the specified control limits, with the exception of the compounds presented in the following table.

Sample Locations	Initial/Continuing	Compound	Criteria
IA-3			
Effluent-1	ICV %RSD	3-Chloropropane	36.3%
Influent-1			

The criteria used to evaluate the initial and continuing calibration are presented in the following table. In the case of a calibration deviation, the sample results are qualified.

Initial/Continuing	Criteria	Sample Result	Qualification
	RRF <0.05	Non-detect	R
	KKF <0.05	Detect	J
Initial and Continuing	RRF <0.01 ¹	Non-detect	R
Calibration	KKF <0.01	Detect	J
	RRF >0.05 or RRF >0.01 ¹	Non-detect	No Action
	KKF >0.05 01 KKF >0.01	Detect	NO ACTION
	%RSD > 15% or a correlation coefficient <0.99	Non-detect	UJ
Initial Calibration	70KSD > 1576 OF a COTTERATION COEMICIENT <0.99	Detect	J
	%RSD >90%	Non-detect	R

Initial/Continuing	Criteria	Sample Result	Qualification
		Detect	J
Continuing Calibration	%D >20% (increase in sensitivity)	Non-detect	No Action
		Detect	J
	%D >20% (decrease in sensitivity)	Non-detect	UJ
		Detect	J
	%D >90% (increase/decrease in sensitivity)	Non-detect	R
		Detect	J

Note:

5. Surrogates/System Monitoring Compounds

All samples to be analyzed for organic compounds are spiked with surrogate compounds prior to sample preparation to evaluate overall laboratory performance and efficiency of the analytical technique. VOC analysis requires that all surrogates associated with the analysis exhibit a percent recovery within the established acceptance limits of 70% to 130%.

Surrogate recoveries were within control limits.

6. Internal Standard Performance

Internal standard performance criteria ensure that the GC/MS sensitivity and response are stable during every sample analysis. The criteria requires the internal standard compounds associated with the VOC exhibit area counts that are not greater than two times (+100%) or less than one-half (-50%) of the area counts of the associated continuing calibration standard.

Internal standard responses were within control limits.

7. Laboratory Control Sample (LCS) Analysis

The LCS analysis is used to assess the accuracy of the analytical method independent of matrix interferences. The compounds associated with the LCS/LCSD analysis must exhibit a percent recovery within the established acceptance limits of 70% to 130%.

Sample locations associated with LCS/LCSD analysis exhibiting recoveries outside of the control limits presented in the following table.

Sample Locations	Compound	LCS Recovery	LCSD Recovery
IA-3			
Effluent-1	3-Chloropropene	>UL	>UL
Influent-1			

¹ RRF of 0.01 only applies to compounds which are typically poor responding compounds (i.e., ketones, 1,4-dioxane, etc.)

The criteria used to evaluate the LCS/LCSD recoveries are presented in the following table. In the case of an LCS/LCSD deviation, the sample results are qualified as documented in the table below.

Control Limit	Sample Result	Qualification
the upper central limit (LIL)	Non-detect	No Action
> the upper control limit (UL)	Detect	J
< the lower control limit (LL) but > 10%	Non-detect	UJ
the lower control limit (LL) but > 10%	Detect	J
< 10%	Non-detect	R
1070	Detect	J

8. Field Duplicate Analysis

Field duplicate analysis is used to assess the overall precision of the field sampling procedures and analytical method. A control limit of 50% for air matrices is applied to the RPD between the parent sample and the field duplicate. In the instance when the parent and/or duplicate sample concentrations are less than or equal to 5 times the RL, a control limit of three times the RL is applied for air matrices.

A field duplicate was not included with this SDG.

9. Compound Identification

Compounds are identified on the GC/MS by using the analytes relative retention time and ion spectra.

Sample results associated with compound that exhibited a concentration greater than the linear range of the instrument calibration are summarized in the following table.

Sample ID	Compound	Original Analysis	Diluted Analysis	Reported Analysis
IA-3	Ethanol	-	1100 E	1100 EJ
Effluent-1	Ethanol	-	840 E	840 EJ
Influent-1	Ethanol	-	560 E	560 EJ

Note: In the instance where both the original analysis and the diluted analysis sample results exhibited a concentration greater than and/or less than the calibration linear range of the instrument; the sample result exhibiting the greatest concentration will be reported as the final result.

Sample results associated with compounds exhibiting concentrations greater than the linear range are qualified as documented in the table below when reported as the final reported sample result.

Reported Sample Results	Qualification
Diluted sample result within calibration range	D

Reported Sample Results	Qualification
Diluted sample result less than the calibration range	DJ
Diluted sample result greater than the calibration range	EDJ
Original sample result greater than the calibration range	EJ

10. System Performance and Overall Assessment

Please note, the laboratory includes a Limit of Detection (LOD) in the laboratory report which is specific to Department of Defense (DOD) reporting and should not be considered for this site/project. Only the Reporting Limit (RL) and Method Detection Limit (MDL) are stored in the database for this data set.

Overall system performance was acceptable. Other than for those deviations specifically mentioned in this review, the overall data quality is within the guidelines specified in the method.

DATA VALIDATION CHECKLIST FOR VOCs

VOCs: TO-15	Re	ported		ormance eptable	Not
	No	Yes	No	Yes	Required
GAS CHROMATOGRAPHY/MASS SPECTROMETE	RY (GC/N	MS)			
Tier II Validation					
Holding times		X		Х	
Canister return pressure (<-1"Hg)		Х		X	
Reporting limits (units)		Х		Х	
Blanks					1
A. Method blanks		Х	Х		
B. Equipment blanks	Х				Х
C. Trip blanks	Х				Х
Laboratory Control Sample (LCS)		Х	Х		
Laboratory Control Sample Duplicate (LCSD)		Х	Х		
LCS/LCSD Precision (RPD)		Х		X	
Matrix Spike (MS)	Х				Х
Matrix Spike Duplicate (MSD)	Х				X
MS/MSD Precision (RPD)	Х				Х
Field/Lab Duplicate (RPD)	Х				Х
Surrogate Spike Recoveries		Х		X	
Dilution Factor		Х		X	
Moisture Content	Х				Х
Tier III Validation					I
System performance and column resolution		X		X	
Initial calibration %RSDs		Х	Х		
Continuing calibration RRFs		X		X	
Continuing calibration %Ds		X		X	
Instrument tune and performance check		X		X	
Ion abundance criteria for each instrument used		X		X	
Internal standard		X		X	
Compound identification and quantitation					
A. Reconstructed ion chromatograms		Х		Х	

VOCs: TO-15	Rep	oorted		rmance ptable	Not
	No	Yes	No	Yes	Required
GAS CHROMATOGRAPHY/MASS SPECTROMETR	Y (GC/M	S)			
B. Quantitation Reports		Х		X	
C. RT of sample compounds within the established RT windows		Х		Х	
D. Transcription/calculation errors present		X		X	
E. Reporting limits adjusted to reflect sample dilutions		Х		Х	

Notes:

%RSD Relative standard deviation

%R Percent recovery

RPD Relative percent difference

%D Percent difference

DATA USABILITY SUMMARY REPORT

SAMPLE COMPLIANCE REPORT

Sample	0					(Compliar	ıcy ¹		
Delivery Group (SDG)	Sampling Date	Protocol	Sample ID	Matrix	voc	svoc	ALD	PFAS	MET	Noncompliance
	5/20/2021	SW846	IA-3	Air	Yes	-	-	-	-	VOC: ICV %RSD
2105446	5/20/2021	SW846	Effluent-1	Air	Yes	-	-	-	-	VOC: ICV %RSD, MB
	5/20/2021	SW846	Influent-1	Air	No	-	-	-	-	VOC: ICV %RSD

Note:

Samples which are compliant with no added validation qualifiers are listed as "yes". Samples which are non-compliant or which have added qualifiers are listed as "no". A "no" designation does not necessarily indicate that the data have been rejected or are otherwise unusable.

DATA USABILITY SUMMARY REPORT

VALIDATION PERFORMED BY: Jeffrey L. Davin

SIGNATURE:

DATE: June 27, 2021

PEER REVIEW: Dennis K. Capria

DATE: June 29, 2021

CHAIN OF CUSTODY CORRECTED SAMPLE ANALYSIS DATA SHEETS



Analysis Request /Canister Chain of Custody

For Laboratory Use Only

Workorder #:

Sample Transportation Notice: Relinquishing signature on this document indicates that samples are shipped in compliance with a any kind. Relinquishing signature also indicates agreement to hold harmless, defend, and indemnify Eurofins Air Toxics against any samples. D.O.T Hotline (800) 467-4922	Shipper Name: Fidey Custody Seals Intact? Yes No		Relinquished by: (Signature/Affiliation) Date Time	Relinquisned by: (Signature/Artiliation)	Arcad13 5/2/21	and by I/Ging through Affiliation					05A +ntluent- 640758 23465 65/20/21	EXT West - 6-0902 23828 * 05/21/21	6LZZ(0 23977 05/20		Lab Field Sample Identification(Location) Can# Flow Informa	Sampler: Don Zuds Site Name: 8 ms Step Name: 8 ms	Project #	Project Name: BMS Building 31/2	
Sample Transportation Notice: Relinquishing signature on this document indicates that samples are shipped in compliance with all applicable local, State, Federal, and international laws, regulations, and ordinances of any kind. Relinquishing signature also indicates agreement to hold harmless, defend, and indemnify Eurofins Air Toxics against any claim, demand, or action, of any kind, related to the collection, handling, of shipping of samples. D.O.T Hotline (800) 467-4922	Is Intact? (Yes No None	Lab Use Only 12.5/2.4		*	1730 Received by: (Signature/Affiliation)						23466	23828 * U5/20/21 15 45 105/20/21 1550	23977 05/20/0829	Date Time Date Time Initial Records Final Re	Flow Information Information at (in psi N2 / 15 in	Lab Use O	Canister Vacuum/Pressure Requested Analyses	Plane Report Result Standard Rush 3/2/ (specify)	Special Instructions/Notes: Turnaround Time (Rush surcharges may apply)





MODIFIED EPA METHOD TO-15 GC/MS SIM/FULL SCAN BMS Building 3VI

Cilent ID: IA-3 Lab ID: 2105446-01A Date/Time Collected: 5/20/21 04:28 PM Media:	IA-3 2105446-01A 5/20/21 04:28 PM	Date/Time Analyzed: Dilution Factor: Instrument/Filename:		5/24/21 06:58 PM 1.42 msd21.i / 21052417		
		MD		Rpt. Limit	Amount	1 [
Compound	CAS#	(£m/gn)	(m/gn)	(sw/gn)	(ng/m3)	
1,2,4-Trichlorobenzene	120-82-1	1.8	2.1	5.3	Not Detected	
1,2,4-Trimethylbenzene	9-63-6	0.13	0.28	0.70	0.19 J	
1,2-Dichlorobenzene	95-50-1	0.11	0.34	0.85	Not Detected	
1,2-Dichloropropane	78-87-5	0.067	0.26	99.0	Not Detected	
1,3,5-Trimethylbenzene	108-67-8	0.15	0.28	0.70	Not Detected	
1,3-Butadiene	106-99-0	0.049	0.12	0.31	Not Detected	
1,3-Dichlorobenzene	541-73-1	0.086	0.34	0.85	Not Detected	

Not Detected

42

1.6 J

3.3

0.20 1.3 0.84

0.074 0.21 0.16 0.16

123-91-1 540-84-1 Not Detected

1.7 2.2 0.70

0.15

107-05-1

67-63-0

622-96-8

108-10-1 67-64-1

4-Methyl-2-pentanone

3-Chloropropene

4-Ethyltoluene

591-78-6

78-93-3

2-Butanone (Methyl Ethyl Ketone)

2-Hexanone 2-Propanol

2,2,4-Trimethylpentane

1,4-Dioxane

100-44-7

75-27-4 75-25-2 74-83-9 75-15-0

Bromodichloromethane

Bromoform

alpha-Chlorotoluene

Acetone

0.12 0.55 0.12 0.10 0.20 0.33 0.39

0.15 J 0.19 J Not Detected
Not Detected
Not Detected
Not Detected
Not Detected
Not Detected

27

0.58 3.4 0.74

1.2 0.70 0.89 0.28 0.23 0.67 0.29 0.38

0.95 1.5 2.8 2.2

0.88

10061-01-5

cis-1,3-Dichloropropene

Cumene

Bromomethane Carbon Disulfide

Chlorobenzene

98-85-8

108-90-7

Not Detected

Page 5 of 37

Not Detected Not Detected





MODIFIED EPA METHOD TO-15 GC/MS SIM/FULL SCAN BMS Building 3VI

1.42 msd21.i / 21052417 5/24/21 06:58 PM Instrument/Filename: Date/Time Analyzed: Dilution Factor: **Lab ID:** 2105446-01A **Date/Time Collected:** 5/20/21 04:28 PM Client ID: Media:

		MDL	ГОР	Rpt. Limit	Amount
Compound	CAS#	(ng/m3)	(ng/m3)	(ng/m3)	(ng/m3)
Cyclohexane	110-82-7	0.13	0.98	2.4	Not Detected
Dibromochloromethane	124-48-1	0.15	0.48	1.2	Not Detected
Ethanol	64-17-5	0.14	0.54	1.3	1100 E ^J
Freon 11	75-69-4	0.070	0.32	0.80	1.4
Freon 113	76-13-1	0.15	0.44	1.1	0.36 J
Heptane	142-82-5	0.18	1.2	2.9	0.60 J
Hexachlorobutadiene	87-68-3	1.4	3.0	7.6	Not Detected
Hexane	110-54-3	0.15	1.0	2.5	0.18 J
Methylene Chloride	75-09-2	0.80	2.0	2.5	Not Detected
Propylbenzene	103-65-1	0.10	0.28	0.70	Not Detected
Styrene	100-42-5	0.061	0.24	09.0	0.14 J
Tetrahydrofuran	109-99-9	0.21	0.84	2.1	0.57 J
trans-1,3-Dichloropropene	10061-02-6	0.068	0.26	0.64	Not Detected

D: Analyte not within the DoD scope of accreditation.

TENTATIVELY IDENTIFIED COMPOUNDS

		Match	TOD	Amount
Compound	CAS#			ppbv
1,2,3-Trimethylbenzene	526-73-8	NA		Not Detected
2-Nitropropane	79-46-9	ΝΑ		Not Detected
4-Chlorotoluene	106-43-4	ΑN		Not Detected
Acetonitrile	75-05-8	ΑN		Not Detected
Benzaldehyde	100-52-7	ΝΑ		Not Detected
bis(2-Chloroethyl) Ether	111-44-4	ΝΑ		Not Detected





MODIFIED EPA METHOD TO-15 GC/MS SIM/FULL SCAN BMS Building 3VI

Date/Time Analyzed: Date/Time Collected: 5/20/21 04:28 PM 2105446-01A Client ID: Lab ID:

6 Liter Summa Canister (SIM Certified)

Media:

msd21.i / 21052417 5/24/21 06:58 PM Instrument/Filename: Dilution Factor:

TENTATIVELY IDENTIFIED COMPOUNDS

		Match	ГОР	Amount
Compound	CAS#			nddd
Cyclohexene	110-83-8	NA		Not Detected
Epichlorohydrin	106-89-8	AN		Not Detected
Ethyl Acetate	141-78-6	AN		Not Detected
Isobutanol	78-83-1	NA		Not Detected
Isopropyl ether	108-20-3	AN		Not Detected
Methacrylonitrile	126-98-7	NA		Not Detected
Methyl Acetate	79-20-9	NA		Not Detected
N,N-Dimethyl Aniline	121-69-7	NA		Not Detected
n-Butanol	71-36-3	AN		Not Detected
Nitrobenzene	98-92-3	NA		Not Detected
Pentane	109-66-0	%98		1.1 NJ
Propylene	115-07-1	Ϋ́		Not Detected
Pyridine	110-86-1	NA		Not Detected

E = Exceeds instrument calibration range.
 J = Estimated value.
 NJ = The identification is based on presumptive evidence; estimated value.
 D: Analyte not within the DoD scope of accreditation.

Surrogates	CAS#	Limits	%Recovery
1,2-Dichloroethane-d4	17060-07-0	70-130	101
4-Bromofluorobenzene	460-00-4	70-130	86
Toluene-d8	2037-26-5	70-130	92





MODIFIED EPA METHOD TO-15 GC/MS SIM/FULL SCAN BMS Building 3VI

Client ID:

Lab ID: 2105446-01B

Date/Time Collected: 5/20/21 04:28 PM

Media: 6 Liter Summa Canister (SIM Certified)

1.42 msd21.i / 21052417sim Dilution Factor: Instrument/Filename:

5/24/21 06:58 PM

Date/Time Analyzed:

Media: 0 Liter Suffiffia	o Liter Suffiffia Caffister (Silvi Certiffed)	Instrument/Filename:		IIISQZ I.1 / Z I USZ4 I / SIIIII		
		MDL	ГОР	Rpt. Limit	Amount	
Compound	CAS#	(ng/m3)	(ng/m3)	(ng/m3)	(ng/m3)	
1,1,1-Trichloroethane	71-55-6	0.0043	0.062	0.15	0.014 J	1
1,1,2,2-Tetrachloroethane	79-34-5	0.025	0.078	0.19	Not Detected	
1,1,2-Trichloroethane	2-00-62	0.014	0.062	0.15	Not Detected	
1,1-Dichloroethane	75-34-3	0.031	0.046	0.11	Not Detected	
1,1-Dichloroethene	75-35-4	0.0041	0.045	0.056	Not Detected	
1,2-Dibromoethane (EDB)	106-93-4	0.014	0.087	0.22	0.024 J	
1,2-Dichloroethane	107-06-2	0.0060	0.046	0.11	0.057 J	
1,4-Dichlorobenzene	106-46-7	0.056	0.068	0.17	Not Detected	
Benzene	71-43-2	0.013	0.036	0.23	0.22 J	
Carbon Tetrachloride	56-23-5	0.0089	0.071	0.18	0.30	
Chloroethane	75-00-3	0.0060	0:030	0.19	0.10 J	
Chloroform	67-66-3	0.011	0.055	0.14	0.24	
Chloromethane	74-87-3	0.18	0.59	1.5	0.98 J	
cis-1,2-Dichloroethene	156-59-2	0.0068	0.045	0.11	0.042 J	
Ethyl Benzene	100-41-4	0.0046	0.049	0.12	0.12	
Freon 114	76-14-2	0.017	0.079	0.20	0.12 J	
Freon 12	75-71-8	0.0052	0.056	3.5	2.4 J	
m,p-Xylene	108-38-3	0.0075	0.049	0.25	0.39	
Methyl tert-butyl ether	1634-04-4	0.0092	0.041	0.51	Not Detected	
Naphthalene	91-20-3	0.094	0.15	0.37	0.61	
o-Xylene	95-47-6	0.0086	0.049	0.12	0.15	
Tetrachloroethene	127-18-4	0.010	0.077	0.19	0.20	
Toluene	108-88-3	0.0098	0.043	0.27	0.52	
trans-1,2-Dichloroethene	156-60-5	0.0058	0.045	0.56	Not Detected	

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MODIFIED EPA METHOD TO-15 GC/MS SIM/FULL SCAN BMS Building 3VI

Client ID:

Lab ID: 2105446-01B **Date/Time Collected:** 5/20/21 04:28 PM

6 Liter Summa Canister (SIM Certified) Media:

5/24/21 06:58 PM Date/Time Analyzed:

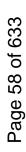
Dilution Factor:

1.42 msd21.i / 21052417sim Instrument/Filename:

		MDL	ГОР	Rpt. Limit	Amount
Compound	CAS#	(ng/m3)	(ng/m3)	(ng/m3)	(ng/m3)
Trichloroethene	79-01-6	0.014	0.061	0.15	0.053 J
Vinyl Chloride	75-01-4	0.0046	0.029	0.036	Not Detected

J = Estimated value. D: Analyte not within the DoD scope of accreditation.

Surrogates	CAS#	Limits	%Recovery	
1,2-Dichloroethane-d4	17060-07-0	70-130	102	
4-Bromofluorobenzene	460-00-4	70-130	88	
Toluene-d8	2037-26-5	70-130	94	





MODIFIED EPA METHOD TO-15 GC/MS SIM/FULL SCAN BMS Building 3VI

1.34 msd21.i / 21052418 5/24/21 07:35 PM Instrument/Filename: Date/Time Analyzed: Dilution Factor: Client ID: Effluent-1
Lab ID: 2105446-02A
Date/Time Collected: 5/20/21 03:50 PM Media:

		MDL	ГОР	Rpt. Limit	Amount
Compound	CAS#	(ng/m3)	(ng/m3)	(ng/m3)	(ng/m3)
1,2,4-Trichlorobenzene	120-82-1	1.7	2.0	5.0	Not Detected
1,2,4-Trimethylbenzene	92-63-6	0.12	0.26	99.0	Not Detected
1,2-Dichlorobenzene	95-50-1	0.11	0.32	0.80	Not Detected
1,2-Dichloropropane	78-87-5	0.063	0.25	0.62	Not Detected
1,3,5-Trimethylbenzene	108-67-8	0.14	0.26	99.0	Not Detected
1,3-Butadiene	106-99-0	0.046	0.12	0.30	Not Detected
1,3-Dichlorobenzene	541-73-1	0.081	0.32	0.80	Not Detected
1,4-Dioxane	123-91-1	0.070	0.19	0.48	0.093 J
2,2,4-Trimethylpentane	540-84-1	0.20	1.2	3.1	Not Detected
2-Butanone (Methyl Ethyl Ketone)	78-93-3	0.15	0.79	2.0	Not Detected
2-Hexanone	591-78-6	0.15	1.	2.7	Not Detected
2-Propanol	67-63-0	0.14	99.0	1.6	88
3-Chloropropene	107-05-1	99.0	0.84	2.1	Not Detected
4-Ethyltoluene	622-96-8	0.11	0.26	99.0	Not Detected
4-Methyl-2-pentanone	108-10-1	0.12	0.22	0.55	Not Detected
Acetone	67-64-1	0.52	0.64	3.2	14
alpha-Chlorotoluene	100-44-7	0.11	0.28	0.69	Not Detected
Bromodichloromethane	75-27-4	0.097	0.36	0.90	Not Detected
Bromoform	75-25-2	0.19	0.55	1.4	Not Detected
Bromomethane	74-83-9	0.31	1.0	2.6	Not Detected
Carbon Disulfide	75-15-0	0.37	0.83	2.1	Not Detected
Chlorobenzene	108-90-7	0.057	0.25	0.62	Not Detected
cis-1,3-Dichloropropene	10061-01-5	090'0	0.24	0.61	Not Detected
Cumene	98-82-8	0.070	0.26	99.0	Not Detected

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MODIFIED EPA METHOD TO-15 GC/MS SIM/FULL SCAN BMS Building 3VI

1.34 msd21.i / 21052418 5/24/21 07:35 PM Instrument/Filename: Date/Time Analyzed: Dilution Factor: Client ID: Effluent-1
Lab ID: 2105446-02A
Date/Time Collected: 5/20/21 03:50 PM Media:

		MDL	ГОР	Rpt. Limit	Amount
Compound	CAS#	(ng/m3)	(ng/m3)	(ng/m3)	(ng/m3)
Cyclohexane	110-82-7	0.12	0.92	2.3	Not Detected
Dibromochloromethane	124-48-1	0.14	0.46	1.1	Not Detected
Ethanol	64-17-5	0.14	0.50	1.3	840 EJ
Freon 11	75-69-4	0.066	0.30	0.75	1.9
Freon 113	76-13-1	0.14	0.41	1.0	0.26 J
Heptane	142-82-5	0.17	1.1	2.7	Not Detected
Hexachlorobutadiene	87-68-3	1.3	2.8	7.1	Not Detected
Hexane	110-54-3	0.14	0.94	2.4	Not Detected
Methylene Chloride	75-09-2	0.75	1.9	2.3	Not Detected
Propylbenzene	103-65-1	0.10	0.26	99.0	Not Detected
Styrene	100-42-5	0.057	0.23	0.57	0.081 J
Tetrahydrofuran	109-99-9	0.20	0.79	2.0	0.36 J
trans-1,3-Dichloropropene	10061-02-6	0.064	0.24	0.61	Not Detected

D: Analyte not within the DoD scope of accreditation.

TENTATIVELY IDENTIFIED COMPOUNDS

		Match	COD	Amount
Compound	CAS#			ppbv
1,2,3-Trimethylbenzene	526-73-8	NA		Not Detected
2-Nitropropane	79-46-9	Ϋ́		Not Detected
4-Chlorotoluene	106-43-4	Ϋ́		Not Detected
Acetonitrile	75-05-8	Ϋ́		Not Detected
Benzaldehyde	100-52-7	Ϋ́		Not Detected
bis(2-Chloroethyl) Ether	111-44-4	٩		Not Detected





MODIFIED EPA METHOD TO-15 GC/MS SIM/FULL SCAN BMS Building 3VI

Client ID: Effluent-1
Lab ID: 2105446-02A
Date/Time Collected: 5/20/21 03:50 PM

Media:

6 Liter Summa Canister (SIM Certified)

5/24/21 07:35 PM Date/Time Analyzed: Dilution Factor:

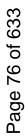
1.34 msd21.i / 21052418 Instrument/Filename:

TENTATIVELY IDENTIFIED COMPOUNDS

		Match	ГОР	Amount
Compound	CAS#			hpbv
Cyclohexene	110-83-8	NA		Not Detected
Epichlorohydrin	106-89-8	Ν		Not Detected
Ethyl Acetate	141-78-6	ΑN		Not Detected
Isobutanol	78-83-1	Ν		Not Detected
Isopropyl ether	108-20-3	Ν		Not Detected
Methacrylonitrile	126-98-7	Ν		Not Detected
Methyl Acetate	79-20-9	Ν		Not Detected
N,N-Dimethyl Aniline	121-69-7	Ν		Not Detected
n-Butanol	71-36-3	ΑN		Not Detected
Nitrobenzene	98-95-3	Ν		Not Detected
Pentane	109-66-0	Ν		Not Detected
Propylene	115-07-1	Ą		Not Detected
Pyridine	110-86-1	Ϋ́		Not Detected

E = Exceeds instrument calibration range.J = Estimated value.D: Analyte not within the DoD scope of accreditation.

Surrogates	CAS#	Limits	%Recovery
1,2-Dichloroethane-d4	17060-07-0	70-130	107
4-Bromofluorobenzene	460-00-4	70-130	87
Toluene-d8	2037-26-5	70-130	92





MODIFIED EPA METHOD TO-15 GC/MS SIM/FULL SCAN BMS Building 3VI

1.34 msd21.i / 21052418sim 5/24/21 07:35 PM Instrument/Filename: Date/Time Analyzed: Dilution Factor: 6 Liter Summa Canister (SIM Certified) Client ID: Effluent-1
Lab ID: 2105446-02B
Date/Time Collected: 5/20/21 03:50 PM Media:

		MDL	ГОР	Rpt. Limit	Amount
Compound	CAS#	(ng/m3)	(ug/m3)	(ng/m3)	(ng/m3)
1,1,1-Trichloroethane	71-55-6	0.0040	0.058	0.15	0.0068 J
1,1,2,2-Tetrachloroethane	79-34-5	0.023	0.074	0.18	Not Detected
1,1,2-Trichloroethane	79-00-5	0.013	0.058	0.15	Not Detected
1,1-Dichloroethane	75-34-3	0.030	0.043	0.11	Not Detected
1,1-Dichloroethene	75-35-4	0.0039	0.042	0.053	Not Detected
1,2-Dibromoethane (EDB)	106-93-4	0.013	0.082	0.20	Not Detected
1,2-Dichloroethane	107-06-2	0.0056	0.043	0.11	U.037 J
1,4-Dichlorobenzene	106-46-7	0.053	0.064	0.16	Not Detected
Benzene	71-43-2	0.012	0.034	0.21	0.21 <u>0.020</u> J UB
Carbon Tetrachloride	56-23-5	0.0084	0.067	0.17	0.092 J
Chloroethane	75-00-3	0.0056	0.028	0.18	0.052 J
Chloroform	67-66-3	0.010	0.052	0.13	0.29
Chloromethane	74-87-3	0.17	0.55	1.4	0.86 J
cis-1,2-Dichloroethene	156-59-2	0.0064	0.042	0.11	0.072 J
Ethyl Benzene	100-41-4	0.0043	0.046	0.12	0.12 0.0048-J UB
Freon 114	76-14-2	0.016	0.075	0.19	0.15 J
Freon 12	75-71-8	0.0049	0.053	3.3	2.2 J
m,p-Xylene	108-38-3	0.0071	0.046	0.23	0.23 0 .023 J- UB
Methyl tert-butyl ether	1634-04-4	0.0087	0.039	0.48	Not Detected
Naphthalene	91-20-3	0.089	0.14	0.35	Not Detected
o-Xylene	95-47-6	0.0081	0.046	0.12	0.013 J
Tetrachloroethene	127-18-4	0.0094	0.073	0.18	0.015 J
Toluene	108-88-3	0.0092	0.040	0.25	0.018 J
trans-1,2-Dichloroethene	156-60-5	0.0055	0.042	0.53	0.013 J

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MODIFIED EPA METHOD TO-15 GC/MS SIM/FULL SCAN BMS Building 3VI

1.34 msd21.i / 21052418sim 5/24/21 07:35 PM Instrument/Filename: Date/Time Analyzed: Dilution Factor: 6 Liter Summa Canister (SIM Certified) Client ID: Effluent-1
Lab ID: 2105446-02B
Date/Time Collected: 5/20/21 03:50 PM Media:

		MDL	ГОР	Rpt. Limit	Amount
Compound	CAS#	(m/gn)	(ng/m3)	(ng/m3)	(ng/m3)
Trichloroethene	79-01-6	0.014	0.058	0.14	Not Detected
Vinyl Chloride	75-01-4	0.0043	0.027	0.034	Not Detected

J = Estimated value. D: Analyte not within the DoD scope of accreditation.

Surrogates	CAS#	Limits	%Recovery
1,2-Dichloroethane-d4	17060-07-0	70-130	101
4-Bromofluorobenzene	460-00-4	70-130	68
Toluene-d8	2037-26-5	70-130	94





MODIFIED EPA METHOD TO-15 GC/MS SIM/FULL SCAN BMS Building 3VI

Client ID: Influent-1
Lab ID: 2105446-03A Date/Time Analyzed: Date/Time Analyzed: Date/Time Collected: 5/20/21 04:15 PM Dilution Factor:

Media:

Dilution Factor: 1.16 Instrument/Filename: msd21.i / 21052419

5/24/21 08:39 PM

		MDL	ГОР	Rpt. Limit	Amount
Compound	CAS#	(ng/m3)	(ng/m3)	(ng/m3)	(ng/m3)
1,2,4-Trichlorobenzene	120-82-1	1.4	1.7	4.3	Not Detected
1,2,4-Trimethylbenzene	95-63-6	0.10	0.23	0.57	0.20 J
1,2-Dichlorobenzene	95-50-1	0.092	0.28	0.70	Not Detected
1,2-Dichloropropane	78-87-5	0.054	0.21	0.54	0.26 J
1,3,5-Trimethylbenzene	108-67-8	0.12	0.23	0.57	Not Detected
1,3-Butadiene	106-99-0	0.040	0.10	0.26	Not Detected
1,3-Dichlorobenzene	541-73-1	0.070	0.28	0.70	Not Detected
1,4-Dioxane	123-91-1	090'0	0.17	0.42	0.14 J
2,2,4-Trimethylpentane	540-84-1	0.17	1.1	2.7	Not Detected
2-Butanone (Methyl Ethyl Ketone)	78-93-3	0.13	0.68	1.7	3.9
2-Hexanone	591-78-6	0.13	0.95	2.4	Not Detected
2-Propanol	67-63-0	0.12	0.57	1.4	49
3-Chloropropene	107-05-1	0.57	0.73	1.8	Not Detected
4-Ethyltoluene	622-96-8	0.096	0.23	0.57	0.17 J
4-Methyl-2-pentanone	108-10-1	0.11	0.19	0.48	0.21 J
Acetone	67-64-1	0.45	0.55	2.8	24
alpha-Chlorotoluene	100-44-7	0.098	0.24	09.0	Not Detected
Bromodichloromethane	75-27-4	0.084	0.31	0.78	Not Detected
Bromoform	75-25-2	0.17	0.48	1.2	Not Detected
Bromomethane	74-83-9	0.27	0.90	2.2	Not Detected
Carbon Disulfide	75-15-0	0.32	0.72	1.8	1.4 J
Chlorobenzene	108-90-7	0.050	0.21	0.53	Not Detected
cis-1,3-Dichloropropene	10061-01-5	0.052	0.21	0.53	Not Detected
Cumene	98-82-8	090'0	0.23	0.57	0.073 J

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MODIFIED EPA METHOD TO-15 GC/MS SIM/FULL SCAN BMS Building 3VI

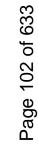
msd21.i / 21052419 5/24/21 08:39 PM Instrument/Filename: Date/Time Analyzed: Dilution Factor: **Lab ID:** 2105446-03A **Date/Time Collected:** 5/20/21 04:15 PM Influent-1 Client ID: Media:

		MDL	ГОР	Rpt. Limit	Amount
Compound	CAS#	(ng/m3)	(ng/m3)	(ng/m3)	(ng/m3)
Cyclohexane	110-82-7	0.10	08'0	2.0	0.36 J
Dibromochloromethane	124-48-1	0.12	0.40	0.99	Not Detected
Ethanol	64-17-5	0.12	0.44	1.1	560 E J
Freon 11	75-69-4	0.057	0.26	0.65	1.4
Freon 113	76-13-1	0.12	0.36	0.89	0.41 J
Heptane	142-82-5	0.15	0.95	2.4	0.50 J
Hexachlorobutadiene	87-68-3	1.1	2.5	6.2	Not Detected
Hexane	110-54-3	0.12	0.82	2.0	0.45 J
Methylene Chloride	75-09-2	0.65	1.6	2.0	0.92 J
Propylbenzene	103-65-1	0.086	0.23	0.57	Not Detected
Styrene	100-42-5	0.050	0.20	0.49	0.078 J
Tetrahydrofuran	109-99-9	0.17	0.68	1.7	0.40 J
trans-1,3-Dichloropropene	10061-02-6	0.055	0.21	0.53	Not Detected

D: Analyte not within the DoD scope of accreditation.

TENTATIVELY IDENTIFIED COMPOUNDS

		Match	ГОР	Amount
Compound	CAS#			ppbv
1,2,3-Trimethylbenzene	526-73-8	NA		Not Detected
1-Propanol, 2-methyl-	78-83-1	20%		1.1 NJ
2-Nitropropane	79-46-9	ΑN		Not Detected
4-Chlorotoluene	106-43-4	ΑN		Not Detected
Acetonitrile	75-05-8	Ϋ́		Not Detected
Benzaldehyde	100-52-7	ΝΑ		Not Detected





MODIFIED EPA METHOD TO-15 GC/MS SIM/FULL SCAN BMS Building 3VI

5/24/21 08:39 PM Date/Time Analyzed: Date/Time Collected: 5/20/21 04:15 PM 2105446-03A Influent-1 Client ID: Lab ID:

6 Liter Summa Canister (SIM Certified)

Media:

msd21.i / 21052419 Instrument/Filename: Dilution Factor:

TENTATIVELY IDENTIFIED COMPOUNDS

		Match	ΓOD	Amount
Compound	CAS#			hpbv
bis(2-Chloroethyl) Ether	111-44-4	NA		Not Detected
Cyclohexene	110-83-8	ΝΑ		Not Detected
Epichlorohydrin	106-89-8	ΝΑ		Not Detected
Ethyl Acetate	141-78-6	ΝΑ		Not Detected
Isopropyl ether	108-20-3	ΑΝ		Not Detected
Methacrylonitrile	126-98-7	ΝΑ		Not Detected
Methyl Acetate	79-20-9	ΝΑ		Not Detected
N,N-Dimethyl Aniline	121-69-7	ΑΝ		Not Detected
n-Butanol	71-36-3	ΝΑ		Not Detected
Nitrobenzene	98-95-3	ΝΑ		Not Detected
Pentane	109-66-0	80%		5.8 NJ
Propylene	115-07-1	ΑΝ		Not Detected
Pyridine	110-86-1	ΝΑ		Not Detected

E = Exceeds instrument calibration range.
J = Estimated value.
NJ = The identification is based on presumptive evidence; estimated value.
D: Analyte not within the DoD scope of accreditation.

Surrogates	CAS#	Limits	%Recovery
1,2-Dichloroethane-d4	17060-07-0	70-130	106
4-Bromofluorobenzene	460-00-4	70-130	81
Toluene-d8	2037-26-5	70-130	94





MODIFIED EPA METHOD TO-15 GC/MS SIM/FULL SCAN BMS Building 3VI

Lab ID: 2105446-03B **Date/Time Collected:** 5/20/21 04:15 PM

Influent-1

Client ID:

6 Liter Summa Canister (SIM Certified) Media:

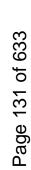
msd21.i / 21052419sim Instrument/Filename: Dilution Factor:

5/24/21 08:39 PM

Date/Time Analyzed:

		MDL	ГОР	Rpt. Limit	Amount
Compound	CAS#	(ng/m3)	(ng/m3)	(ng/m3)	(ng/m3)
1,1,1-Trichloroethane	71-55-6	0.0035	0.051	0.13	0.014 J
1,1,2,2-Tetrachloroethane	79-34-5	0.020	0.064	0.16	Not Detected
1,1,2-Trichloroethane	2-00-62	0.012	0.051	0.13	Not Detected
1,1-Dichloroethane	75-34-3	0.026	0.038	0.094	Not Detected
1,1-Dichloroethene	75-35-4	0.0034	0.037	0.046	Not Detected
1,2-Dibromoethane (EDB)	106-93-4	0.012	0.071	0.18	0.021 J
1,2-Dichloroethane	107-06-2	0.0049	0.038	0.094	U.087 J
1,4-Dichlorobenzene	106-46-7	0.046	0.056	0.14	Not Detected
Benzene	71-43-2	0.010	0:030	0.18	0.23
Carbon Tetrachloride	56-23-5	0.0073	0.058	0.14	0.32
Chloroethane	75-00-3	0.0049	0.024	0.15	0.051 J
Chloroform	67-66-3	0.0089	0.045	0.11	0.36
Chloromethane	74-87-3	0.14	0.48	1.2	0.93 J
cis-1,2-Dichloroethene	156-59-2	0.0056	0.037	0.092	0.039 J
Ethyl Benzene	100-41-4	0.0037	0.040	0.10	0.19
Freon 114	76-14-2	0.014	0.065	0.16	0.12 J
Freon 12	75-71-8	0.0042	0.046	2.9	2.3 J
m,p-Xylene	108-38-3	0.0061	0.040	0.20	0.52
Methyl tert-butyl ether	1634-04-4	0.0075	0.033	0.42	0.032 J
Naphthalene	91-20-3	0.077	0.12	0:30	0.67
o-Xylene	95-47-6	0.0070	0.040	0.10	0.22
Tetrachloroethene	127-18-4	0.0082	0.063	0.16	0.18
Toluene	108-88-3	0.0080	0.035	0.22	16
trans-1,2-Dichloroethene	156-60-5	0.0048	0.037	0.46	0.012 J

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💸 eurofins

BMS Building 3VI

msd21.i / 21052419sim 5/24/21 08:39 PM Instrument/Filename: Date/Time Analyzed: Dilution Factor: 6 Liter Summa Canister (SIM Certified) **Lab ID:** 2105446-03B **Date/Time Collected:** 5/20/21 04:15 PM Influent-1 Client ID: Media:

#		בפ	Rpt. Limit	Amonnt
3 00 02	CAS# (ug/m3)	(ng/m3)	(ng/m3)	(ng/m3)
0-10-67	79-01-6 0.012	0.050	0.12	0.058 J
Vinyl Chloride 75-01-4 0	0.0037	0.024	0.030	Not Detected

J = Estimated value. D: Analyte not within the DoD scope of accreditation.

Surrogates	CAS#	Limits	%Recovery
1,2-Dichloroethane-d4	17060-07-0	70-130	103
4-Bromofluorobenzene	460-00-4	70-130	84
Toluene-d8	2037-26-5	70-130	93



Bristol Myers Squibb Thompson Road Investigation

Data Usability Summary Report

Syracuse, NY

Volatile Organic Compound (VOC) Analysis

SDGs # 2101359

Analyses Performed By: Eurofins Air Toxics

Folsom, CA

Report #40371R Review Level: Tier III Project: 30064943

SUMMARY

This data quality assessment summarizes the review of Sample Delivery Group (SDG) # 2101359 for samples collected in association with the Bristol Myers Squibb Thompson Road Site. The review was conducted as a Tier III evaluation and included review of data package completeness. Only analytical data associated with constituents of concern were reviewed for this validation. Field documentation was not included in this review. Included with this assessment are the validation annotated sample result sheets, and chain of custody. Analyses were performed on the following samples:

0			Sample	Parent			Analysis	;	
Sample ID	Lab ID	Matrix	Collection Date	Sample	voc	svoc	PFAS	MET	ALD
IA-3	2011359-01A	Air	1/14/2021		Х				
AMB-011421	2011359-02A	Air	1/14/2021		Х				
DUP-011421	2011359-03A	Air	1/14/2021	IA-3	Х				

ANALYTICAL DATA PACKAGE DOCUMENTATION

The table below is the evaluation of the data package completeness.

	Rep	orted		rmance ptable	Not
Items Reviewed	No	Yes	No	Yes	Required
Sample receipt condition		Х		Х	
2. Requested analyses and sample results		Х		Х	
Master tracking list		Х		Х	
4. Methods of analysis		Х		Х	
5. Reporting limits		Х		Х	
6. Sample collection date		Х		Х	
7. Laboratory sample received date		Х		Х	
8. Sample preservation verification (as applicable)		Х		Х	
Sample preparation/extraction/analysis dates		Х		Х	
10. Fully executed Chain-of-Custody (COC) form		Х		X	
11. Narrative summary of QA or sample problems provided		Х		X	
12. Data Package Completeness and Compliance		Х		Х	

Note:

QA - Quality Assurance

List of Acronyms

%D: Percent Difference

%R: Percent Recovery

AC: Acceptable

ALC/GLY: Alcohols/Glycols

BAL: Blank Action Level

CCV: Continuing Calibration Verification

CRDL: Contract Required Detection Limit

D: Dilution

EIS: Extractable Internal Standard

FB: Field Blank

FD: Field Duplicate

ALD: Aldehydes

GC/ECD: Gas Chromatograph/Electron Capture Detector

GC/MS: Gas Chromatograph/Mass spectrometer

HT: Holding Time

ICP: Inductively Coupled Plasma

ICS: Interference Control Sample

ICV: Initial Calibration Verification

ISTD: Internal Standards

LabDup: Laboratory Duplicate

LCS: Lab Control Sample

LCSD: Lab Control Sample Duplicate

LL: Lower Control Limit

MB: Method Blank

MDL: Method Detection Limit

MET: Metals

MS: Matrix Spike

MSD: Matrix Spike Duplicate

N/A: Not Applicable

NC: Not Compliant

List of Acronyms, Continued

PAH: Polyaromatic Hydrocarbon

PCB: Polychlorinated Biphenyl

PEST: Pesticide

PFAS: Per- and Polyfluoroalkyl Substances

QA: Quality Assurance

QC: Quality Control

RB: Rinse Blank

RL: Reporting Limit

RPD: Relative Percent Difference

RRF: Relative Response Factor

RSD: Relative Standard Deviation

RT: Retention Time

SDG: Sample Delivery Group

SerDil: Serial Dilution

SIM: Single Ion Monitoring

SOP: Standard Operating Procedure

SSTD: Surrogate Standards

SVOC: Semivolatile Organic Compound

TB: Trip Blank

TIC: Tentatively Identified Compound

TOC: Total Organic Carbon

TOTDIS: Total and Dissolved

UL: Upper Control Limit

USEPA: United States Environmental Protection Agency

VOC: Volatile Organic Compound

ORGANIC ANALYSIS INTRODUCTION

Analyses were performed according to United States Environmental Protection Agency (USEPA) Method TO-15. Data were reviewed in accordance with USEPA National Functional Guidelines of October 1999, USEPA Region II SOP HW-31- Validating Air Samples Volatile Organic Analysis of Ambient Air In Canister by Method TO-15 of October 2006, New York State DEC Analytical Method ASP 2005 TO-15 (QA/QC Criteria R9 TO-15), and NYSDEC Modifications to R9 TO-15 QA/QC Criteria October 2009.

The data review process is an evaluation of data on a technical basis rather than a determination of contract compliance. As such, the standards against which the data are being weighed may differ from those specified in the analytical method. It is assumed that the data package represents the best efforts of the laboratory and had already been subjected to adequate and sufficient quality review prior to submission.

During the review process, laboratory qualified and unqualified data are verified against the supporting documentation. Based on this evaluation, qualifier codes may be added, deleted, or modified by the data reviewer. Results are qualified with the following codes in accordance with USEPA National Functional Guidelines:

- Concentration (C) Qualifiers
 - U The compound was analyzed for but not detected. The associated value is the compound quantitation limit.
 - B The compound has been found in the sample as well as its associated blank, its presence in the sample may be suspect.
- Quantitation (Q) Qualifiers
 - E The compound was quantitated above the calibration range.
 - D Concentration is based on a diluted sample analysis.
- Validation Qualifiers
 - J The analyte was positively identified; however, the associated numerical value is an estimated concentration only.
 - UJ The analyte was not detected above the reported sample detection limit. However, the reported limit is approximate and may or may not represent the actual limit of detection.
 - UB Analyte considered non-detect at the listed value due to associated blank contamination.
 - N The analysis indicates the presence of a compound for which there is presumptive evidence to make a tentative identification.
 - R The sample results are rejected.

Two facts should be noted by all data users. First, the "R" flag means that the associated value is unusable. In other words, due to significant quality control (QC) problems, the analysis is invalid and provides no information as to whether the compound is present or not. "R" values should not appear on data tables because they cannot be relied upon, even as a last resort. The second fact to keep in mind is that no compound concentration, even if it has passed all QC tests, is guaranteed to be accurate. Strict QC serves to increase confidence in data but any value potentially contains error.

VOLATILE ORGANIC COMPOUND (VOC) ANALYSES

1. Holding Times

The specified holding times for the following methods are presented in the following table.

Method	Matrix	Holding Time	Preservation	Return Canister Pressure
USEPA TO-15	Air	30 days from collection to analysis	Ambient Temperature	< -1" Hg

All samples were analyzed within the specified holding time and canister return pressure / vacuum criteria.

2. Blank Contamination

Quality assurance (QA) blanks (i.e., method and rinse blanks) are prepared to identify any contamination which may have been introduced into the samples during sample preparation or field activity. Method blanks measure laboratory contamination. Rinse blanks measure contamination of samples during field operations.

A blank action level (BAL) of five times the concentration of a detected compound in an associated blank (common laboratory contaminant compounds are calculated at ten times) is calculated for QA blanks containing concentrations greater than the method detection limit (MDL). The BAL is compared to the associated sample results to determine the appropriate qualification of the sample results, if needed.

All compounds associated with the QA blanks exhibited a concentration less than the MDL, with the exception of the compounds listed in the following table. Sample results associated with QA blank contamination that were greater than the BAL resulted in the removal of the laboratory qualifier (B) of data. Sample results less than the BAL associated with the following sample locations were qualified as listed in the following table.

Sample Locations	Analytes	Sample Result	Qualification
IA-3	Methylene chloride	Detected sample results <rl <bal<="" and="" td=""><td>"UB" at the RL</td></rl>	"UB" at the RL
AMB-011421	2-Propanol		

Note:

RL Reporting limit

3. Mass Spectrometer Tuning

Mass spectrometer performance was acceptable and all analyses were performed within a 12-hour tune clock.

System performance and column resolution were acceptable.

4. Calibration

Satisfactory instrument calibration is established to ensure that the instrument is capable of producing acceptable quantitative data. An initial calibration demonstrates that the instrument is capable of acceptable performance at the beginning of an experimental sequence. The continuing calibration verifies that the instrument daily performance is satisfactory.

4.1 Initial Calibration

The method specifies percent relative standard deviation (%RSD) and relative response factor (RRF) limits for select compounds only. A technical review of the data applies limits to all compounds with no exceptions.

All target compounds associated with the initial calibration standards must exhibit a %RSD less than the control limit (30%) or a correlation coefficient greater than 0.99 and an RRF value greater than control limit (0.05).

4.2 Continuing Calibration

All target compounds associated with the continuing calibration standard must exhibit a percent difference (%D) less than the control limit (30%) and RRF value greater than control limit (0.05).

Compounds associated with the calibrations were within the specified control limits, with the exception of the compounds presented in the following table.

Sample Locations	Initial/Continuing	Compound	Criteria
IA-3			
AMB-011421	ICV %RSD	3-Chloropropane	35.2%
DUP-011421			

The criteria used to evaluate the initial and continuing calibration are presented in the following table. In the case of a calibration deviation, the sample results are qualified.

Initial/Continuing	Criteria	Sample Result	Qualification	
	RRF <0.05	Non-detect	R	
	KKF <0.05	Detect	J	
Initial and Continuing Calibration	RRF <0.01 ¹	Non-detect	R	
	KKF <0.01	Detect	J	
	RRF >0.05 or RRF >0.01 ¹	Non-detect	No Action	
	KKF >0.05 01 KKF >0.01	Detect	NO ACTION	
	%RSD > 15% or a correlation coefficient <0.99	Non-detect	UJ	
Initial Calibration	%RSD > 15% or a correlation coefficient <0.99	Detect	J	
	0/DCD - 000/	Non-detect	R	
	%RSD >90%	Detect	J	

Initial/Continuing	Criteria	Sample Result	Qualification
	%D >20% (increase in sensitivity)	Non-detect	No Action
Continuing Calibration	70D >20 /0 (IIIClease III sellsluvity)	Detect	J
	%D >20% (decrease in sensitivity)	Non-detect	UJ
	70D >20 /0 (decrease in sensitivity)	Detect	J
	9/D > 009/ (increase/degreese in consitivity)	Non-detect	R
	%D >90% (increase/decrease in sensitivity)	Detect	J

Note:

5. Surrogates/System Monitoring Compounds

All samples to be analyzed for organic compounds are spiked with surrogate compounds prior to sample preparation to evaluate overall laboratory performance and efficiency of the analytical technique. VOC analysis requires that all surrogates associated with the analysis exhibit a percent recovery within the established acceptance limits of 70% to 130%.

Surrogate recoveries were within control limits.

Internal Standard Performance

Internal standard performance criteria ensure that the GC/MS sensitivity and response are stable during every sample analysis. The criteria requires the internal standard compounds associated with the VOC exhibit area counts that are not greater than two times (+100%) or less than one-half (-50%) of the area counts of the associated continuing calibration standard.

Internal standard responses were within control limits.

7. Laboratory Control Sample (LCS) Analysis

The LCS analysis is used to assess the accuracy of the analytical method independent of matrix interferences. The compounds associated with the LCS/LCSD analysis must exhibit a percent recovery within the established acceptance limits of 70% to 130%.

Compounds associated with the LCS analysis exhibited recoveries within the control limits.

8. Field Duplicate Analysis

Field duplicate analysis is used to assess the overall precision of the field sampling procedures and analytical method. A control limit of 50% for air matrices is applied to the RPD between the parent sample and the field duplicate. In the instance when the parent and/or duplicate sample concentrations are less than or equal to 5 times the RL, a control limit of three times the RL is applied for air matrices.

Results for duplicate samples are summarized in the following table (ug/m3).

		Sample	Duplicate	
Sample ID/Duplicate ID	Analyte	Result	Result	RPD
	1,2,4-Trimethylbenzene	0.16 J	0.17 J	AC

¹ RRF of 0.01 only applies to compounds which are typically poor responding compounds (i.e., ketones, 1,4-dioxane, etc.)

		Sample	Duplicate	
Sample ID/Duplicate ID	Analyte	Result	Result	RPD
	1,4-Dioxane	0.065 J	0.13 J	AC
	2-Butanone (Methyl Ethyl Ketone)	0.94 J	1.3 J	AC
	2-Propanol	19	18	5.4%
	4-Ethyltoluene	0.14 J	0.14 J	AC
	Acetone	7.9	7.6	AC
	Bromodichloromethane	0.057 J	0.061 J	AC
	Cumene	0.053 J	0.057 J	AC
	Ethanol	580 E	550 E	5.3%
	Freon 11	1.4	1.2	AC
	Freon 113	0.52 J	0.47 J	AC
	Heptane	0.29 J	0.27 J	AC
	Hexane	0.54 J	0.42 J	AC
	Styrene	0.045 J	0.053 J	AC
	Pentane	1.6	1.3	20.7
	1,1,1-Trichloroethane	0.32 J	0.32 J	AC
-3/	1,2-Dibromoethane (EDB)	0.027 J	0.22 U	AC
JP-011421	1,2-Dichloroethane	0.069 J	0.068 J	AC
	Benzene	0.67	0.63	AC
	Carbon Tetrachloride	0.44	0.42	AC
	Chloroethane	0.041 J	0.046 J	AC
	Chloroform	0.38	0.38	AC
	Chloromethane	0.64 J	0.64 J	AC
	cis-1,2-Dichloroethene	0.069 J	0.064 J	AC
	Ethyl Benzene	0.11 J	0.10 J	AC
	Freon 114	0.10 J	0.098 J	AC
	Freon 12	2.3 J	2.3 J	AC
	m,p-Xylene	0.35	0.30	AC
	Naphthalene	0.22 J	0.26 J	AC
	o-Xylene	0.12	0.12 J	AC
	Tetrachloroethene	0.74	0.26	AC
	Toluene	0.67	0.65	AC
	Trichloroethene	0.54	0.52	AC

Notes:

AC = Acceptable

NC = Not Compliant

The calculated RPDs between the parent sample and field duplicate were acceptable.

9. Compound Identification

Compounds are identified on the GC/MS by using the analytes relative retention time and ion spectra.

Sample results associated with compound that exhibited a concentration greater than the linear range of the instrument calibration are summarized in the following table.

Sample ID	Compound	Original Analysis	Diluted Analysis	Reported Analysis
IA-3	Ethanol	-	580 E	580 EJ
DUP-011421	Ethanol	-	550 E	550 EJ

Note: In the instance where both the original analysis and the diluted analysis sample results exhibited a concentration greater than and/or less than the calibration linear range of the instrument; the sample result exhibiting the greatest concentration will be reported as the final result.

Sample results associated with compounds exhibiting concentrations greater than the linear range are qualified as documented in the table below when reported as the final reported sample result.

Reported Sample Results	Qualification
Diluted sample result within calibration range	D
Diluted sample result less than the calibration range	DJ
Diluted sample result greater than the calibration range	EDJ
Original sample result greater than the calibration range	EJ

10. System Performance and Overall Assessment

Please note, the laboratory includes a Limit of Detection (LOD) in the laboratory report which is specific to Department of Defense (DOD) reporting and should not be considered for this site/project. Only the Reporting Limit (RL) and Method Detection Limit (MDL) are stored in the database for this data set.

Overall system performance was acceptable. Other than for those deviations specifically mentioned in this review, the overall data quality is within the guidelines specified in the method.

DATA VALIDATION CHECKLIST FOR VOCs

VOCs: TO-15		Reported		Performance Acceptable	
		Yes	No	Yes	Required
GAS CHROMATOGRAPHY/MASS SPECTROM	ETRY (GC/N	/IS)			
Tier II Validation					
Holding times		X		X	
Canister return pressure (<-1"Hg)		Х		X	
Reporting limits (units)		Х		X	
Blanks					
A. Method blanks		Х	Х		
B. Equipment blanks	Х				Х
C. Trip blanks	Х				Х
Laboratory Control Sample (LCS)		Х		Х	
Laboratory Control Sample Duplicate (LCSD)	Х				Х
LCS/LCSD Precision (RPD)	Х				Х
Matrix Spike (MS)	Х				Х
Matrix Spike Duplicate (MSD)	Х				Х
MS/MSD Precision (RPD)	Х				Х
Field/Lab Duplicate (RPD)		Х		X	
Surrogate Spike Recoveries		Х		Х	
Dilution Factor		X		X	
Moisture Content		Х		X	
Tier III Validation	ı			1	I
System performance and column resolution		X		X	
Initial calibration %RSDs		Х	Х		
Continuing calibration RRFs		X		Х	
Continuing calibration %Ds		Х		X	
Instrument tune and performance check		X		X	
Ion abundance criteria for each instrument used		X		Х	
Internal standard		X		Х	
Compound identification and quantitation					
A. Reconstructed ion chromatograms		X		X	

VOCs: TO-15		Reported		rmance ptable	Not
		Yes	No	Yes	Required
GAS CHROMATOGRAPHY/MASS SPECTROMETRY (GC/MS)					
B. Quantitation Reports		Х		X	
C. RT of sample compounds within the established RT windows		Х		Х	
D. Transcription/calculation errors present		X		X	
E. Reporting limits adjusted to reflect sample dilutions		Х		Х	

Notes:

%RSD Relative standard deviation

%R Percent recovery

RPD Relative percent difference

%D Percent difference

DATA USABILITY SUMMARY REPORT

SAMPLE COMPLIANCE REPORT

Sample	0					(Compliar	ncy ¹		
Delivery Group (SDG)	Sampling Date	Protocol	Sample ID	Matrix	voc	svoc	ALD	PFAS	MET	Noncompliance
	1/14/2021	SW846	IA-3	Air	Yes	No	Yes	-	-	VOC: MB, ICV %RSD
2101359	1/14/2021	SW846	AMB-011421	Air	Yes	No	Yes	-	-	VOC: MB, ICV %RSD
	1/14/2021	SW846	DUP-011421	Air	No	No	Yes	-	-	VOC: ICV %RSD

Note:

Samples which are compliant with no added validation qualifiers are listed as "yes". Samples which are non-compliant or which have added qualifiers are listed as "no". A "no" designation does not necessarily indicate that the data have been rejected or are otherwise unusable.

DATA USABILITY SUMMARY REPORT

VALIDATION PERFORMED BY: Jeffrey L. Davin

SIGNATURE:

DATE: February 28, 2021

PEER REVIEW: Dennis K. Capria

DATE: March 3, 2021

CHAIN OF CUSTODY CORRECTED SAMPLE ANALYSIS DATA SHEETS



Analysis Request /Canister Chain of Custody

For Laboratory Use Only

180 Blue Ravine Rd. Suite B, Folsom, CA 95630 Phone (800) 985-5955; Fax (916) 351-8279	330 PID:	Workorder #:	# 1365	537	2101359	9		an-		
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Shipper Name: Wake How	Custody Seals Intact?	(Yes	No	None	3-15 15					
sample Transportation Notice: Relinquishing signature on this document indicates that samples are shipped in compliance with all any kind. Relinquishing signature also indicates agreement to hold harmless, defend, and indemnify European Air Taxing parties any kind. Relinquishing signature also indicates agreement to hold harmless, defend, and indemnify European Air Taxing parties are	ature on this document in ement to hold harmless	ndicates that sam	iples are shipped	d in complianc		le local, State	, Federal, and	applicable local, State, Federal, and international laws, regulations, and ordinances of	gulations, and ordin	ances of
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* See went order for List of Companiels



Client ID: IA-3

Lab ID: 2101359-01A **Date/Time Analyzed:** 1/25/21 09:27 PM

Date/Time Collected: 1/14/21 04:42 PM Dilution Factor: 1.42

Media: Instrument/Filename: msd21.i / 21012523

		MDL	LOD	Rpt. Limit	Amount
Compound	CAS#	(ug/m3)	(ug/m3)	(ug/m3)	(ug/m3)
1,2,4-Trichlorobenzene	120-82-1	1.2	1.6	5.3	Not Detected
1,2,4-Trimethylbenzene	95-63-6	0.049	0.17	0.70	0.16 J
1,2-Dichlorobenzene	95-50-1	0.098	0.21	0.85	Not Detected
1,2-Dichloropropane	78-87-5	0.036	0.16	0.66	Not Detected
1,3,5-Trimethylbenzene	108-67-8	0.047	0.17	0.70	Not Detected
1,3-Butadiene	106-99-0	0.017	0.078	0.31	Not Detected
1,3-Dichlorobenzene	541-73-1	0.060	0.21	0.85	Not Detected
1,4-Dioxane	123-91-1	0.035	0.13	0.51	0.065 J
2,2,4-Trimethylpentane	540-84-1	0.35	1.0	3.3	Not Detected
2-Butanone (Methyl Ethyl Ketone)	78-93-3	0.15	0.63	2.1	0.94 J
2-Hexanone	591-78-6	0.20	0.87	2.9	Not Detected
2-Propanol	67-63-0	0.30	0.52	1.7	19
3-Chloropropene	107-05-1	0.53	0.67	4.4	Not Detected J
4-Ethyltoluene	622-96-8	0.035	0.17	0.70	0.14 J
4-Methyl-2-pentanone	108-10-1	0.054	0.14	0.58	Not Detected
Acetone	67-64-1	0.43	0.50	3.4	7.9
alpha-Chlorotoluene	100-44-7	0.13	0.18	0.74	Not Detected
Bromodichloromethane	75-27-4	0.029	0.24	0.95	0.057 J
Bromoform	75-25-2	0.094	0.37	1.5	Not Detected
Bromomethane	74-83-9	0.36	0.83	2.8	Not Detected
Carbon Disulfide	75-15-0	0.39	0.66	2.2	Not Detected
Chlorobenzene	108-90-7	0.020	0.16	0.65	Not Detected
cis-1,3-Dichloropropene	10061-01-5	0.031	0.16	0.64	Not Detected
Cumene	98-82-8	0.024	0.17	0.70	0.053 J



Client ID: IA-3

Lab ID: 2101359-01A **Date/Time Analyzed:** 1/25/21 09:27 PM

Date/Time Collected: 1/14/21 04:42 PM Dilution Factor: 1.42

Media: Instrument/Filename: msd21.i / 21012523

		MDL	LOD	Rpt. Limit	Amount
Compound	CAS#	(ug/m3)	(ug/m3)	(ug/m3)	(ug/m3)
Cyclohexane	110-82-7	0.26	0.73	2.4	Not Detected
Dibromochloromethane	124-48-1	0.078	0.30	1.2	Not Detected
Ethanol	64-17-5	0.31	0.40	1.3	580 E J
Freon 11	75-69-4	0.033	0.20	0.80	1.4
Freon 113	76-13-1	0.11	0.27	1.1	0.52 J
Heptane	142-82-5	0.098	0.87	2.9	0.29 J
Hexachlorobutadiene	87-68-3	1.6	2.3	7.6	Not Detected
Hexane	110-54-3	0.19	0.75	2.5	0.54 J
Methylene Chloride	75-09-2	0.49	0.74	0.99	0.99 0 .50 J UB
Propylbenzene	103-65-1	0.044	0.17	0.70	Not Detected
Styrene	100-42-5	0.024	0.15	0.60	0.045 J
Tetrahydrofuran	109-99-9	0.22	0.63	2.1	Not Detected
trans-1,3-Dichloropropene	10061-02-6	0.039	0.16	0.64	Not Detected

D: Analyte not within the DoD scope of accreditation.

TENTATIVELY IDENTIFIED COMPOUNDS

		Match	LOD	Amount	
Compound	CAS#			ppbv	
1,2,3-Trimethylbenzene	526-73-8	NA		Not Detected	
2-Nitropropane	79-46-9	NA		Not Detected	
4-Chlorotoluene	106-43-4	NA		Not Detected	
Acetonitrile	75-05-8	NA		Not Detected	
Benzaldehyde	100-52-7	NA		Not Detected	
bis(2-Chloroethyl) Ether	111-44-4	NA		Not Detected	



Client ID: IA-3

Lab ID: 2101359-01A **Date/Time Analyzed:** 1/25/21 09:27 PM

Date/Time Collected: 1/14/21 04:42 PM Dilution Factor: 1.42

Media: 6 Liter Summa Canister (SIM Certified) Instrument/Filename: msd21.i / 21012523

TENTATIVELY IDENTIFIED COMPOUNDS

		Match	LOD	Amount	
Compound	CAS#			ppbv	
Cyclohexene	110-83-8	NA		Not Detected	
Epichlorohydrin	106-89-8	NA		Not Detected	
Ethyl Acetate	141-78-6	NA		Not Detected	
Isobutanol	78-83-1	NA		Not Detected	
Isopropyl ether	108-20-3	NA		Not Detected	
Methacrylonitrile	126-98-7	NA		Not Detected	
Methyl Acetate	79-20-9	NA		Not Detected	
N,N-Dimethyl Aniline	121-69-7	NA		Not Detected	
n-Butanol	71-36-3	NA		Not Detected	
Nitrobenzene	98-95-3	NA		Not Detected	
Pentane	109-66-0	72%		1.6	
Propylene	115-07-1	NA		Not Detected	
Pyridine	110-86-1	NA		Not Detected	

E = Exceeds instrument calibration range.

D: Analyte not within the DoD scope of accreditation.

Surrogates	CAS#	Limits	%Recovery
1,2-Dichloroethane-d4	17060-07-0	70-130	111
4-Bromofluorobenzene	460-00-4	70-130	82
Toluene-d8	2037-26-5	70-130	96

J = Estimated value.



Client ID: IA-3

Lab ID: 2101359-01B **Date/Time Analyzed:** 1/25/21 09:27 PM

Date/Time Collected: 1/14/21 04:42 PM Dilution Factor: 1.42

Media: 6 Liter Summa Canister (SIM Certified) Instrument/Filename: msd21.i / 21012523sim

		MDL	LOD	Rpt. Limit	Amount
Compound	CAS#	(ug/m3)	(ug/m3)	(ug/m3)	(ug/m3)
1,1,1-Trichloroethane	71-55-6	0.0092	0.062	0.15	0.032 J
1,1,2,2-Tetrachloroethane	79-34-5	0.038	0.078	0.19	Not Detected
1,1,2-Trichloroethane	79-00-5	0.017	0.062	0.15	Not Detected
1,1-Dichloroethane	75-34-3	0.033	0.046	0.11	Not Detected
1,1-Dichloroethene	75-35-4	0.019	0.045	0.056	Not Detected
1,2-Dibromoethane (EDB)	106-93-4	0.026	0.087	0.22	0.027 J
1,2-Dichloroethane	107-06-2	0.017	0.046	0.11	0.069 J
1,4-Dichlorobenzene	106-46-7	0.085	0.13	0.17	Not Detected
Benzene	71-43-2	0.022	0.036	0.23	0.67
Carbon Tetrachloride	56-23-5	0.066	0.071	0.18	0.44
Chloroethane	75-00-3	0.012	0.030	0.19	0.041 J
Chloroform	67-66-3	0.022	0.055	0.14	0.38
Chloromethane	74-87-3	0.019	0.023	1.5	0.64 J
cis-1,2-Dichloroethene	156-59-2	0.020	0.045	0.11	0.069 J
Ethyl Benzene	100-41-4	0.019	0.049	0.12	0.11 J
Freon 114	76-14-2	0.013	0.079	0.20	0.10 J
Freon 12	75-71-8	0.016	0.056	3.5	2.3 J
m,p-Xylene	108-38-3	0.025	0.049	0.25	0.35
Methyl tert-butyl ether	1634-04-4	0.022	0.041	0.51	Not Detected
Naphthalene	91-20-3	0.15	0.28	0.37	0.22 J
o-Xylene	95-47-6	0.026	0.049	0.12	0.12
Tetrachloroethene	127-18-4	0.0077	0.077	0.19	0.74
Toluene	108-88-3	0.017	0.043	0.27	0.67
trans-1,2-Dichloroethene	156-60-5	0.017	0.045	0.56	Not Detected



Client ID: IA-3

Lab ID: 2101359-01B **Date/Time Analyzed:** 1/25/21 09:27 PM

Date/Time Collected: 1/14/21 04:42 PM Dilution Factor: 1.42

Media: 6 Liter Summa Canister (SIM Certified) Instrument/Filename: msd21.i / 21012523sim

Compound	CAS#	MDL (ug/m3)	LOD (ug/m3)	Rpt. Limit (ug/m3)	Amount (ug/m3)
Trichloroethene	79-01-6	0.015	0.061	0.15	0.54
Vinyl Chloride	75-01-4	0.0083	0.029	0.036	Not Detected

J = Estimated value.

D: Analyte not within the DoD scope of accreditation.

Surrogates	CAS#	Limits	%Recovery
1,2-Dichloroethane-d4	17060-07-0	70-130	94
4-Bromofluorobenzene	460-00-4	70-130	92
Toluene-d8	2037-26-5	70-130	92



 Client ID:
 AMB-011421

 Lab ID:
 2101359-02A

Date/Time Collected: 1/14/21 04:40 PM

Media:

Date/Time Analyzed: 1/25/21 08:14 PM

Dilution Factor: 1.40

Instrument/Filename: msd21.i / 21012521

		MDL	LOD	Rpt. Limit	Amount
Compound	CAS#	(ug/m3)	(ug/m3)	(ug/m3)	(ug/m3)
1,2,4-Trichlorobenzene	120-82-1	1.2	1.6	5.2	Not Detected
1,2,4-Trimethylbenzene	95-63-6	0.048	0.17	0.69	0.15 J
1,2-Dichlorobenzene	95-50-1	0.097	0.21	0.84	Not Detected
1,2-Dichloropropane	78-87-5	0.036	0.16	0.65	Not Detected
1,3,5-Trimethylbenzene	108-67-8	0.047	0.17	0.69	Not Detected
1,3-Butadiene	106-99-0	0.017	0.077	0.31	Not Detected
1,3-Dichlorobenzene	541-73-1	0.059	0.21	0.84	Not Detected
1,4-Dioxane	123-91-1	0.034	0.13	0.50	0.047 J
2,2,4-Trimethylpentane	540-84-1	0.34	0.98	3.3	Not Detected
2-Butanone (Methyl Ethyl Ketone)	78-93-3	0.14	0.62	2.1	0.50 J
2-Hexanone	591-78-6	0.20	0.86	2.9	Not Detected
2-Propanol	67-63-0	0.30	0.52	1.7	1.7 -1.4 J UB
3-Chloropropene	107-05-1	0.52	0.66	4.4	Not Detected J
4-Ethyltoluene	622-96-8	0.035	0.17	0.69	0.15 J
4-Methyl-2-pentanone	108-10-1	0.054	0.14	0.57	Not Detected
Acetone	67-64-1	0.42	0.50	3.3	4.1
alpha-Chlorotoluene	100-44-7	0.13	0.18	0.72	Not Detected
Bromodichloromethane	75-27-4	0.028	0.23	0.94	Not Detected
Bromoform	75-25-2	0.093	0.36	1.4	Not Detected
Bromomethane	74-83-9	0.36	0.82	2.7	Not Detected
Carbon Disulfide	75-15-0	0.39	0.65	2.2	Not Detected
Chlorobenzene	108-90-7	0.020	0.16	0.64	Not Detected
cis-1,3-Dichloropropene	10061-01-5	0.030	0.16	0.64	Not Detected
Cumene	98-82-8	0.024	0.17	0.69	Not Detected



 Client ID:
 AMB-011421

 Lab ID:
 2101359-02A

Date/Time Collected: 1/14/21 04:40 PM

Media:

Date/Time Analyzed: 1/25/21 08:14 PM

Dilution Factor: 1.40

Instrument/Filename: msd21.i / 21012521

		MDL	LOD	Rpt. Limit	Amount
Compound	CAS#	(ug/m3)	(ug/m3)	(ug/m3)	(ug/m3)
Cyclohexane	110-82-7	0.25	0.72	2.4	Not Detected
Dibromochloromethane	124-48-1	0.077	0.30	1.2	Not Detected
Ethanol	64-17-5	0.31	0.40	1.3	3.4
Freon 11	75-69-4	0.032	0.20	0.79	1.2
Freon 113	76-13-1	0.11	0.27	1.1	0.43 J
Heptane	142-82-5	0.097	0.86	2.9	0.23 J
Hexachlorobutadiene	87-68-3	1.6	2.2	7.5	Not Detected
Hexane	110-54-3	0.19	0.74	2.5	0.36 J
Methylene Chloride	75-09-2	0.49	0.73	0.97	Not Detected
Propylbenzene	103-65-1	0.044	0.17	0.69	Not Detected
Styrene	100-42-5	0.024	0.15	0.60	0.027 J
Tetrahydrofuran	109-99-9	0.22	0.62	2.1	Not Detected
trans-1,3-Dichloropropene	10061-02-6	0.039	0.16	0.64	Not Detected

D: Analyte not within the DoD scope of accreditation.

TENTATIVELY IDENTIFIED COMPOUNDS

	_	Match	LOD	Amount	
Compound	CAS#			ppbv	
1,2,3-Trimethylbenzene	526-73-8	NA		Not Detected	
2-Nitropropane	79-46-9	NA		Not Detected	
4-Chlorotoluene	106-43-4	NA		Not Detected	
Acetonitrile	75-05-8	NA		Not Detected	
Benzaldehyde	100-52-7	NA		Not Detected	
bis(2-Chloroethyl) Ether	111-44-4	NA		Not Detected	



Client ID: AMB-011421 **Lab ID:** 2101359-02A

Date/Time Collected: 1/14/21 04:40 PM

Media: 6 Liter Summa Canister (SIM Certified)

Date/Time Analyzed: 1/25/21 08:14 PM

Dilution Factor: 1.40

Instrument/Filename: msd21.i / 21012521

TENTATIVELY IDENTIFIED COMPOUNDS

		Match	LOD	Amount	
Compound	CAS#			ppbv	
Cyclohexene	110-83-8	NA		Not Detected	
Epichlorohydrin	106-89-8	NA		Not Detected	
Ethyl Acetate	141-78-6	NA		Not Detected	
Isobutanol	78-83-1	NA		Not Detected	
Isopropyl ether	108-20-3	NA		Not Detected	
Methacrylonitrile	126-98-7	NA		Not Detected	
Methyl Acetate	79-20-9	NA		Not Detected	
N,N-Dimethyl Aniline	121-69-7	NA		Not Detected	
n-Butanol	71-36-3	NA		Not Detected	
Nitrobenzene	98-95-3	NA		Not Detected	
Pentane	109-66-0	NA		Not Detected	
Propylene	115-07-1	NA		Not Detected	
Pyridine	110-86-1	NA		Not Detected	

J = Estimated value.

D: Analyte not within the DoD scope of accreditation.

Surrogates	CAS#	Limits	%Recovery
1,2-Dichloroethane-d4	17060-07-0	70-130	104
4-Bromofluorobenzene	460-00-4	70-130	84
Toluene-d8	2037-26-5	70-130	93



Client ID: AMB-011421 **Lab ID:** 2101359-02B

Date/Time Collected: 1/14/21 04:40 PM

Media: 6 Liter Summa Canister (SIM Certified)

Date/Time Analyzed: 1/25/21 08:14 PM

Dilution Factor: 1.40

Instrument/Filename: msd21.i / 21012521sim

		MDL	LOD	Rpt. Limit	Amount
Compound	CAS#	(ug/m3)	(ug/m3)	(ug/m3)	(ug/m3)
1,1,1-Trichloroethane	71-55-6	0.0091	0.061	0.15	Not Detected
1,1,2,2-Tetrachloroethane	79-34-5	0.038	0.077	0.19	Not Detected
1,1,2-Trichloroethane	79-00-5	0.016	0.061	0.15	Not Detected
1,1-Dichloroethane	75-34-3	0.033	0.045	0.11	Not Detected
1,1-Dichloroethene	75-35-4	0.018	0.044	0.056	Not Detected
1,2-Dibromoethane (EDB)	106-93-4	0.026	0.086	0.22	Not Detected
1,2-Dichloroethane	107-06-2	0.017	0.045	0.11	0.10 J
1,4-Dichlorobenzene	106-46-7	0.084	0.13	0.17	Not Detected
Benzene	71-43-2	0.022	0.036	0.22	0.65
Carbon Tetrachloride	56-23-5	0.066	0.070	0.18	0.46
Chloroethane	75-00-3	0.012	0.030	0.18	0.027 J
Chloroform	67-66-3	0.022	0.055	0.14	0.070 J
Chloromethane	74-87-3	0.019	0.023	1.4	0.62 J
cis-1,2-Dichloroethene	156-59-2	0.019	0.044	0.11	Not Detected
Ethyl Benzene	100-41-4	0.018	0.049	0.12	0.10 J
Freon 114	76-14-2	0.012	0.078	0.20	0.10 J
Freon 12	75-71-8	0.016	0.055	3.5	2.3 J
m,p-Xylene	108-38-3	0.025	0.049	0.24	0.29
Methyl tert-butyl ether	1634-04-4	0.022	0.040	0.50	Not Detected
Naphthalene	91-20-3	0.15	0.28	0.37	Not Detected
o-Xylene	95-47-6	0.025	0.049	0.12	0.11 J
Tetrachloroethene	127-18-4	0.0076	0.076	0.19	0.13 J
Toluene	108-88-3	0.017	0.042	0.26	0.65
trans-1,2-Dichloroethene	156-60-5	0.016	0.044	0.56	Not Detected



Client ID: AMB-011421 **Lab ID:** 2101359-02B

Date/Time Collected: 1/14/21 04:40 PM

Media: 6 Liter Summa Canister (SIM Certified)

Date/Time Analyzed: 1/25/21 08:14 PM

Dilution Factor: 1.40

Instrument/Filename: msd21.i / 21012521sim

		MDL	LOD	Rpt. Limit	Amount
Compound	CAS#	(ug/m3)	(ug/m3)	(ug/m3)	(ug/m3)
Trichloroethene	79-01-6	0.015	0.060	0.15	0.023 J
Vinyl Chloride	75-01-4	0.0082	0.029	0.036	Not Detected

J = Estimated value.

D: Analyte not within the DoD scope of accreditation.

Surrogates	CAS#	Limits	%Recovery
1,2-Dichloroethane-d4	17060-07-0	70-130	95
4-Bromofluorobenzene	460-00-4	70-130	92
Toluene-d8	2037-26-5	70-130	91



 Client ID:
 DUP-011421

 Lab ID:
 2101359-03A

Date/Time Collected: 1/14/21 12:00 AM

Media:

Date/Time Analyzed: 1/25/21 08:50 PM

Dilution Factor: 1.46

Instrument/Filename: msd21.i / 21012522

		MDL	LOD	Rpt. Limit	Amount
Compound	CAS#	(ug/m3)	(ug/m3)	(ug/m3)	(ug/m3)
1,2,4-Trichlorobenzene	120-82-1	1.2	1.6	5.4	Not Detected
1,2,4-Trimethylbenzene	95-63-6	0.050	0.18	0.72	0.17 J
1,2-Dichlorobenzene	95-50-1	0.10	0.22	0.88	Not Detected
1,2-Dichloropropane	78-87-5	0.038	0.17	0.67	Not Detected
1,3,5-Trimethylbenzene	108-67-8	0.049	0.18	0.72	Not Detected
1,3-Butadiene	106-99-0	0.018	0.081	0.32	Not Detected
1,3-Dichlorobenzene	541-73-1	0.061	0.22	0.88	Not Detected
1,4-Dioxane	123-91-1	0.036	0.13	0.53	0.13 J
2,2,4-Trimethylpentane	540-84-1	0.36	1.0	3.4	Not Detected
2-Butanone (Methyl Ethyl Ketone)	78-93-3	0.15	0.64	2.2	1.3 J
2-Hexanone	591-78-6	0.21	0.90	3.0	Not Detected
2-Propanol	67-63-0	0.31	0.54	1.8	18
3-Chloropropene	107-05-1	0.54	0.68	4.6	Not Detected J
4-Ethyltoluene	622-96-8	0.036	0.18	0.72	0.14 J
4-Methyl-2-pentanone	108-10-1	0.056	0.15	0.60	Not Detected
Acetone	67-64-1	0.44	0.52	3.5	7.6
alpha-Chlorotoluene	100-44-7	0.14	0.19	0.76	Not Detected
Bromodichloromethane	75-27-4	0.030	0.24	0.98	0.061 J
Bromoform	75-25-2	0.097	0.38	1.5	Not Detected
Bromomethane	74-83-9	0.37	0.85	2.8	Not Detected
Carbon Disulfide	75-15-0	0.40	0.68	2.3	Not Detected
Chlorobenzene	108-90-7	0.020	0.17	0.67	Not Detected
cis-1,3-Dichloropropene	10061-01-5	0.032	0.16	0.66	Not Detected
Cumene	98-82-8	0.025	0.18	0.72	0.057 J



Client ID: DUP-011421 **Lab ID:** 2101359-03A

Date/Time Collected: 1/14/21 12:00 AM

Media:

Date/Time Analyzed: 1/25/21 08:50 PM

Dilution Factor: 1.46

Instrument/Filename: msd21.i / 21012522

		MDL	LOD	Rpt. Limit	Amount
Compound	CAS#	(ug/m3)	(ug/m3)	(ug/m3)	(ug/m3)
Cyclohexane	110-82-7	0.26	0.75	2.5	Not Detected
Dibromochloromethane	124-48-1	0.080	0.31	1.2	Not Detected
Ethanol	64-17-5	0.32	0.41	1.4	550 E J
Freon 11	75-69-4	0.034	0.20	0.82	1.2
Freon 113	76-13-1	0.12	0.28	1.1	0.47 J
Heptane	142-82-5	0.10	0.90	3.0	0.27 J
Hexachlorobutadiene	87-68-3	1.7	2.3	7.8	Not Detected
Hexane	110-54-3	0.20	0.77	2.6	0.42 J
Methylene Chloride	75-09-2	0.51	0.76	1.0	Not Detected
Propylbenzene	103-65-1	0.046	0.18	0.72	Not Detected
Styrene	100-42-5	0.025	0.16	0.62	0.053 J
Tetrahydrofuran	109-99-9	0.23	0.64	2.2	Not Detected
trans-1,3-Dichloropropene	10061-02-6	0.040	0.16	0.66	Not Detected

D: Analyte not within the DoD scope of accreditation.

TENTATIVELY IDENTIFIED COMPOUNDS

		Match	LOD	Amount	
Compound	CAS#			ppbv	
1,2,3-Trimethylbenzene	526-73-8	NA		Not Detected	
2-Nitropropane	79-46-9	NA		Not Detected	
4-Chlorotoluene	106-43-4	NA		Not Detected	
Acetonitrile	75-05-8	NA		Not Detected	
Benzaldehyde	100-52-7	NA		Not Detected	
bis(2-Chloroethyl) Ether	111-44-4	NA		Not Detected	



Client ID: DUP-011421 **Lab ID:** 2101359-03A

Date/Time Collected: 1/14/21 12:00 AM

Media: 6 Liter Summa Canister (SIM Certified)

Date/Time Analyzed: 1/25/21 08:50 PM

Dilution Factor: 1.46

Instrument/Filename: msd21.i / 21012522

TENTATIVELY IDENTIFIED COMPOUNDS

		Match	LOD	Amount	
Compound	CAS#			ppbv	
Cyclohexene	110-83-8	NA		Not Detected	
Epichlorohydrin	106-89-8	NA		Not Detected	
Ethyl Acetate	141-78-6	NA		Not Detected	
Isobutanol	78-83-1	NA		Not Detected	
Isopropyl ether	108-20-3	NA		Not Detected	
Methacrylonitrile	126-98-7	NA		Not Detected	
Methyl Acetate	79-20-9	NA		Not Detected	
N,N-Dimethyl Aniline	121-69-7	NA		Not Detected	
n-Butanol	71-36-3	NA		Not Detected	
Nitrobenzene	98-95-3	NA		Not Detected	
Pentane	109-66-0	78%		1.3	
Propylene	115-07-1	NA		Not Detected	
Pyridine	110-86-1	NA		Not Detected	

E = Exceeds instrument calibration range.

D: Analyte not within the DoD scope of accreditation.

Surrogates	CAS#	Limits	%Recovery
1,2-Dichloroethane-d4	17060-07-0	70-130	105
4-Bromofluorobenzene	460-00-4	70-130	84
Toluene-d8	2037-26-5	70-130	94

J = Estimated value.



Client ID: DUP-011421 **Lab ID:** 2101359-03B

Date/Time Collected: 1/14/21 12:00 AM

Media: 6 Liter Summa Canister (SIM Certified)

Date/Time Analyzed: 1/25/21 08:50 PM

Dilution Factor: 1.46

Instrument/Filename: msd21.i / 21012522sim

		MDL	LOD	Rpt. Limit	Amount
Compound	CAS#	(ug/m3)	(ug/m3)	(ug/m3)	(ug/m3)
1,1,1-Trichloroethane	71-55-6	0.0095	0.064	0.16	0.032 J
1,1,2,2-Tetrachloroethane	79-34-5	0.039	0.080	0.20	Not Detected
1,1,2-Trichloroethane	79-00-5	0.017	0.064	0.16	Not Detected
1,1-Dichloroethane	75-34-3	0.034	0.047	0.12	Not Detected
1,1-Dichloroethene	75-35-4	0.019	0.046	0.058	Not Detected
1,2-Dibromoethane (EDB)	106-93-4	0.027	0.090	0.22	Not Detected
1,2-Dichloroethane	107-06-2	0.018	0.047	0.12	0.068 J
1,4-Dichlorobenzene	106-46-7	0.087	0.13	0.18	Not Detected
Benzene	71-43-2	0.023	0.037	0.23	0.63
Carbon Tetrachloride	56-23-5	0.068	0.073	0.18	0.42
Chloroethane	75-00-3	0.012	0.031	0.19	0.046 J
Chloroform	67-66-3	0.023	0.057	0.14	0.38
Chloromethane	74-87-3	0.020	0.024	1.5	0.64 J
cis-1,2-Dichloroethene	156-59-2	0.020	0.046	0.12	0.064 J
Ethyl Benzene	100-41-4	0.019	0.051	0.13	0.10 J
Freon 114	76-14-2	0.013	0.082	0.20	0.098 J
Freon 12	75-71-8	0.017	0.058	3.6	2.3 J
m,p-Xylene	108-38-3	0.026	0.051	0.25	0.30
Methyl tert-butyl ether	1634-04-4	0.023	0.042	0.53	Not Detected
Naphthalene	91-20-3	0.16	0.29	0.38	0.26 J
o-Xylene	95-47-6	0.026	0.051	0.13	0.12 J
Tetrachloroethene	127-18-4	0.0079	0.079	0.20	0.26
Toluene	108-88-3	0.018	0.044	0.28	0.65
trans-1,2-Dichloroethene	156-60-5	0.017	0.046	0.58	Not Detected



Client ID: DUP-011421 **Lab ID:** 2101359-03B

Date/Time Collected: 1/14/21 12:00 AM

Media: 6 Liter Summa Canister (SIM Certified)

Date/Time Analyzed: 1/25/21 08:50 PM

Dilution Factor: 1.46

Instrument/Filename: msd21.i / 21012522sim

		MDL	LOD	Rpt. Limit	Amount
Compound	CAS#	(ug/m3)	(ug/m3)	(ug/m3)	(ug/m3)
Trichloroethene	79-01-6	0.016	0.063	0.16	0.52
Vinyl Chloride	75-01-4	0.0086	0.030	0.037	Not Detected

J = Estimated value.

D: Analyte not within the DoD scope of accreditation.

Surrogates	CAS#	Limits	%Recovery
1,2-Dichloroethane-d4	17060-07-0	70-130	93
4-Bromofluorobenzene	460-00-4	70-130	90
Toluene-d8	2037-26-5	70-130	92



Bristol Myers Squibb Thompson Road Investigation

Data Usability Summary Report

Syracuse, NY

Volatile Organic Compound (VOC) Analysis

SDGs # 2106711

Analyses Performed By: Eurofins Air Toxics Folsom, CA

Report #42166R Review Level: Tier III Project: 30064943

SUMMARY

This data quality assessment summarizes the review of Sample Delivery Group (SDG) # 2106711 for samples collected in association with the Bristol Myers Squibb Thompson Road Site. The review was conducted as a Tier III evaluation and included review of data package completeness. Only analytical data associated with constituents of concern were reviewed for this validation. Field documentation was not included in this review. Included with this assessment are the validation annotated sample result sheets, and chain of custody. Analyses were performed on the following samples:

0			Sample	Sample Parent	Analysis				
Sample ID	Lab ID	Matrix			voc	svoc	PFAS	MET	ALD
IA-3	2106711-01A	Air	6/24/2021	DUP-062421	X				
AA-3	2106711-02A	Air	6/24/2021		X				
DUP-062421	2106711-03A	Air	6/24/2021		Х				

ANALYTICAL DATA PACKAGE DOCUMENTATION

The table below is the evaluation of the data package completeness.

	Rep	Reported		rmance ptable	Not
Items Reviewed	No	Yes	No	Yes	Required
Sample receipt condition		Х		Х	
2. Requested analyses and sample results		Х		Х	
Master tracking list		Х		Х	
4. Methods of analysis		Х		Х	
5. Reporting limits		Х		Х	
6. Sample collection date		Х		Х	
7. Laboratory sample received date		Х		Х	
8. Sample preservation verification (as applicable)		Х		Х	
9. Sample preparation/extraction/analysis dates		Х		Х	
10. Fully executed Chain-of-Custody (COC) form		Х		Х	
11. Narrative summary of QA or sample problems provided		Х		X	
12. Data Package Completeness and Compliance		Х		Х	

Note:

QA - Quality Assurance

List of Acronyms

%D: Percent Difference

%R: Percent Recovery

AC: Acceptable

ALC/GLY: Alcohols/Glycols

BAL: Blank Action Level

CCV: Continuing Calibration Verification

CRDL: Contract Required Detection Limit

D: Dilution

EIS: Extractable Internal Standard

FB: Field Blank

FD: Field Duplicate

ALD: Aldehydes

GC/ECD: Gas Chromatograph/Electron Capture Detector

GC/MS: Gas Chromatograph/Mass spectrometer

HT: Holding Time

ICP: Inductively Coupled Plasma

ICS: Interference Control Sample

ICV: Initial Calibration Verification

ISTD: Internal Standards

LabDup: Laboratory Duplicate

LCS: Lab Control Sample

LCSD: Lab Control Sample Duplicate

LL: Lower Control Limit

MB: Method Blank

MDL: Method Detection Limit

MET: Metals

MS: Matrix Spike

MSD: Matrix Spike Duplicate

N/A: Not Applicable

NC: Not Compliant

List of Acronyms, Continued

PAH: Polyaromatic Hydrocarbon

PCB: Polychlorinated Biphenyl

PEST: Pesticide

PFAS: Per- and Polyfluoroalkyl Substances

QA: Quality Assurance

QC: Quality Control

RB: Rinse Blank

RL: Reporting Limit

RPD: Relative Percent Difference

RRF: Relative Response Factor

RSD: Relative Standard Deviation

RT: Retention Time

SDG: Sample Delivery Group

SerDil: Serial Dilution

SIM: Single Ion Monitoring

SOP: Standard Operating Procedure

SSTD: Surrogate Standards

SVOC: Semivolatile Organic Compound

TB: Trip Blank

TIC: Tentatively Identified Compound

TOC: Total Organic Carbon

TOTDIS: Total and Dissolved

UL: Upper Control Limit

USEPA: United States Environmental Protection Agency

VOC: Volatile Organic Compound

ORGANIC ANALYSIS INTRODUCTION

Analyses were performed according to United States Environmental Protection Agency (USEPA) Method TO-15. Data were reviewed in accordance with USEPA National Functional Guidelines of October 1999, USEPA Region II SOP HW-31- Validating Air Samples Volatile Organic Analysis of Ambient Air In Canister by Method TO-15 of October 2006, New York State DEC Analytical Method ASP 2005 TO-15 (QA/QC Criteria R9 TO-15), and NYSDEC Modifications to R9 TO-15 QA/QC Criteria October 2009.

The data review process is an evaluation of data on a technical basis rather than a determination of contract compliance. As such, the standards against which the data are being weighed may differ from those specified in the analytical method. It is assumed that the data package represents the best efforts of the laboratory and had already been subjected to adequate and sufficient quality review prior to submission.

During the review process, laboratory qualified and unqualified data are verified against the supporting documentation. Based on this evaluation, qualifier codes may be added, deleted, or modified by the data reviewer. Results are qualified with the following codes in accordance with USEPA National Functional Guidelines:

- Concentration (C) Qualifiers
 - U The compound was analyzed for but not detected. The associated value is the compound quantitation limit.
 - B The compound has been found in the sample as well as its associated blank, its presence in the sample may be suspect.
- Quantitation (Q) Qualifiers
 - E The compound was quantitated above the calibration range.
 - D Concentration is based on a diluted sample analysis.
- Validation Qualifiers
 - J The analyte was positively identified; however, the associated numerical value is an estimated concentration only.
 - UJ The analyte was not detected above the reported sample detection limit. However, the reported limit is approximate and may or may not represent the actual limit of detection.
 - UB Analyte considered non-detect at the listed value due to associated blank contamination.
 - N The analysis indicates the presence of a compound for which there is presumptive evidence to make a tentative identification.
 - R The sample results are rejected.

Two facts should be noted by all data users. First, the "R" flag means that the associated value is unusable. In other words, due to significant quality control (QC) problems, the analysis is invalid and provides no information as to whether the compound is present or not. "R" values should not appear on data tables because they cannot be relied upon, even as a last resort. The second fact to keep in mind is that no compound concentration, even if it has passed all QC tests, is guaranteed to be accurate. Strict QC serves to increase confidence in data but any value potentially contains error.

VOLATILE ORGANIC COMPOUND (VOC) ANALYSES

1. Holding Times

The specified holding times for the following methods are presented in the following table.

Method	Matrix	Holding Time	Preservation	Return Canister Pressure
USEPA TO-15	Air	30 days from collection to analysis	Ambient Temperature	< -1" Hg

All samples were analyzed within the specified holding time and canister return pressure / vacuum criteria.

2. Blank Contamination

Quality assurance (QA) blanks (i.e., method and rinse blanks) are prepared to identify any contamination which may have been introduced into the samples during sample preparation or field activity. Method blanks measure laboratory contamination. Rinse blanks measure contamination of samples during field operations.

A blank action level (BAL) of five times the concentration of a detected compound in an associated blank (common laboratory contaminant compounds are calculated at ten times) is calculated for QA blanks containing concentrations greater than the method detection limit (MDL). The BAL is compared to the associated sample results to determine the appropriate qualification of the sample results, if needed.

Compounds were not detected above the MDL in the associated blanks; therefore, detected sample results were not associated with blank contamination.

3. Mass Spectrometer Tuning

Mass spectrometer performance was acceptable and all analyses were performed within a 12-hour tune clock.

System performance and column resolution were acceptable.

4. Calibration

Satisfactory instrument calibration is established to ensure that the instrument is capable of producing acceptable quantitative data. An initial calibration demonstrates that the instrument is capable of acceptable performance at the beginning of an experimental sequence. The continuing calibration verifies that the instrument daily performance is satisfactory.

4.1 Initial Calibration

The method specifies percent relative standard deviation (%RSD) and relative response factor (RRF) limits for select compounds only. A technical review of the data applies limits to all compounds with no exceptions.

All target compounds associated with the initial calibration standards must exhibit a %RSD less than the control limit (30%) or a correlation coefficient greater than 0.99 and an RRF value greater than control limit (0.05).

4.2 Continuing Calibration

All target compounds associated with the continuing calibration standard must exhibit a percent difference (%D) less than the control limit (30%) and RRF value greater than control limit (0.05).

Compounds associated with the calibrations were within the specified control limits, with the exception of the compounds presented in the following table.

Sample Locations	Initial/Continuing	Compound	Criteria
IA-3	ICV %RSD	3-Chloropropane	36.3%
AA-3			
DUP-062421	CCV %D	3-Chloropropane	33.0%

The criteria used to evaluate the initial and continuing calibration are presented in the following table. In the case of a calibration deviation, the sample results are qualified.

Initial/Continuing	Criteria	Sample Result	Qualification
	RRF <0.05	Non-detect	R
	NN	Detect	J
Initial and Continuing	RRF <0.01 ¹	Non-detect	R
Calibration	KKI 30.01	Detect	J
	RRF >0.05 or RRF >0.01 ¹	Non-detect	No Action
	KKF >0.05 01 KKF >0.01	Detect	NO ACTION
	%RSD > 15% or a correlation coefficient <0.99	Non-detect	UJ
Initial Calibration	70K3D > 13 % of a confeation coefficient <0.99	Detect	J
ITIIIIAI CAIIDIAIIOII	%RSD >90%	Non-detect	R
	/MGD >90 //0	Detect	J
	9/D > 209/ (increase in constitutiv)	Non-detect	No Action
	%D >20% (increase in sensitivity)	Detect	J
Continuing Calibration	9/D - 209/ (degraded in conditivity)	Non-detect	UJ
Continuing Calibration	%D >20% (decrease in sensitivity)	Detect	J
	9/D > 009/ (increase/degreese in consistivity)	Non-detect	R
	%D >90% (increase/decrease in sensitivity)	Detect	J

Note:

5. Surrogates/System Monitoring Compounds

¹ RRF of 0.01 only applies to compounds which are typically poor responding compounds (i.e., ketones, 1,4-dioxane, etc.)

All samples to be analyzed for organic compounds are spiked with surrogate compounds prior to sample preparation to evaluate overall laboratory performance and efficiency of the analytical technique. VOC analysis requires that all surrogates associated with the analysis exhibit a percent recovery within the established acceptance limits of 70% to 130%.

Surrogate recoveries were within control limits.

6. Internal Standard Performance

Internal standard performance criteria ensure that the GC/MS sensitivity and response are stable during every sample analysis. The criteria requires the internal standard compounds associated with the VOC exhibit area counts that are not greater than two times (+100%) or less than one-half (-50%) of the area counts of the associated continuing calibration standard.

Internal standard responses were within control limits.

7. Laboratory Control Sample (LCS) Analysis

The LCS analysis is used to assess the accuracy of the analytical method independent of matrix interferences. The compounds associated with the LCS/LCSD analysis must exhibit a percent recovery within the established acceptance limits of 70% to 130%.

Sample locations associated with LCS/LCSD analysis exhibiting recoveries outside of the control limits presented in the following table.

Sample Locations	Compound	LCS Recovery	LCSD Recovery	
IA-3	3-Chloropropene	>UL	>UL	
AA-3		111 / 400/	11.1.4.4007	
DUP-062421 Ethanol		<ll but="">10%</ll>	<ll but="">10%</ll>	

The criteria used to evaluate the LCS/LCSD recoveries are presented in the following table. In the case of an LCS/LCSD deviation, the sample results are qualified as documented in the table below.

Control Limit	Sample Result	Qualification
> the upper control limit (UL)	Non-detect	No Action
> the upper control little (OL)	Detect	J
the lower central limit /LL) but > 109/	Non-detect	UJ
< the lower control limit (LL) but > 10%	Detect	J
< 10%	Non-detect	R
< 1076	Detect	J

8. Field Duplicate Analysis

Field duplicate analysis is used to assess the overall precision of the field sampling procedures and analytical method. A control limit of 50% for air matrices is applied to the RPD between the parent

sample and the field duplicate. In the instance when the parent and/or duplicate sample concentrations are less than or equal to 5 times the RL, a control limit of three times the RL is applied for air matrices.

Results for duplicate samples are summarized in the following table.

Sample ID/Duplicate ID	Compound	Sample Result	Duplicate Result	RPD
	1,4-Dioxane	0.14 J	0.19 J	AC
	2-Butanone	1.1 J	1.4 J	AC
	2-Hexanone	2.9 U	0.19 J	AC
	2-Propanol	6.2	6.4	AC
	4-Methyl-2-pentanone	0.14 J	0.15 J	AC
	Acetone	11	12	AC
	Carbon Disulfide	2.2 U	0.65 J	AC
	Cyclohexane	2.4 U	1.6 J	AC
	Ethanol	100	100	0%
	Freon 11	1.1	1.1	AC
	Freon 113	0.42 J	0.38 J	AC
	Styrene	0.070 J	0.093 J	AC
	Tetrahydrofuran	2.1 U	0.22 J	AC
IA-3/DUP-062421	1,1,1-Trichloroethane	0.013 J	0.015 J	AC
	1,2-Dichloroethane	0.042 J	0.043 J	AC
	Benzene	0.11 J	0.12 J	AC
	Carbon Tetrachloride	0.32	0.33	AC
	Chloroethane	0.041 J	0.026 J	AC
	Chloroform	0.14	0.15	AC
	Chloromethane	0.74 J	0.74 J	AC
	cis-1,2-Dichloroethene	0.021 J	0.022 J	AC
	Ethyl Benzene	0.058 J	0.065 J	AC
	Freon 114	0.092 J	0.095 J	AC
	Freon 12	2.2 J	2.2 J	AC
	m,p-Xylene	0.20 J	0.22 J	AC
	Naphthalene	0.43	0.42	AC
	o-Xylene	0.071 J	0.093 J	AC
	Tetrachloroethene	0.082 J	0.060 J	AC

Sample ID/Duplicate ID	Compound	Sample Result	Duplicate Result	RPD
	Toluene	0.45	0.51	AC
	Trichloroethene	0.16	0.15	AC

The calculated RPDs between the parent sample and field duplicate were acceptable.

9. Compound Identification

Compounds are identified on the GC/MS by using the analytes relative retention time and ion spectra. All identified compounds met the specified criteria.

10. System Performance and Overall Assessment

Please note, the laboratory includes a Limit of Detection (LOD) in the laboratory report which is specific to Department of Defense (DOD) reporting and should not be considered for this site/project. Only the Reporting Limit (RL) and Method Detection Limit (MDL) are stored in the database for this data set.

Overall system performance was acceptable. Other than for those deviations specifically mentioned in this review, the overall data quality is within the guidelines specified in the method.

DATA VALIDATION CHECKLIST FOR VOCs

VOCs: TO-15	Re	ported		ormance eptable	Not
	No	Yes	No	Yes	Required
GAS CHROMATOGRAPHY/MASS SPECTROMETE	RY (GC/N	MS)			
Tier II Validation					
Holding times		X		Х	
Canister return pressure (<-1"Hg)		Х		Х	
Reporting limits (units)		Х		Х	
Blanks		'			
A. Method blanks		Х		Х	
B. Equipment blanks	Х				Х
C. Trip blanks	Х				Х
Laboratory Control Sample (LCS)		Х	Х		
Laboratory Control Sample Duplicate (LCSD)		Х	Х		
LCS/LCSD Precision (RPD)		Х		Х	
Matrix Spike (MS)	Х				Х
Matrix Spike Duplicate (MSD)	Х				Х
MS/MSD Precision (RPD)	Х				Х
Field/Lab Duplicate (RPD)	Х				Х
Surrogate Spike Recoveries		Х		X	
Dilution Factor		Х		X	
Moisture Content	Х				Х
Tier III Validation					I
System performance and column resolution		X		X	
Initial calibration %RSDs		Х	Х		
Continuing calibration RRFs		X		X	
Continuing calibration %Ds		X		X	
Instrument tune and performance check		X		X	
Ion abundance criteria for each instrument used		X		X	
Internal standard		X		X	
Compound identification and quantitation					
A. Reconstructed ion chromatograms		Х		X	

VOCs: TO-15	Rep	Reported Performan Acceptab			Not
	No	Yes	No	Yes	Required
GAS CHROMATOGRAPHY/MASS SPECTROMETR	Y (GC/M	S)			
B. Quantitation Reports		X		Х	
C. RT of sample compounds within the established RT windows		Х		Х	
D. Transcription/calculation errors present		X		X	
E. Reporting limits adjusted to reflect sample dilutions		Х		Х	

Notes:

%RSD Relative standard deviation

%R Percent recovery

RPD Relative percent difference

%D Percent difference

DATA USABILITY SUMMARY REPORT

SAMPLE COMPLIANCE REPORT

Sample	0					(Compliar	ıcy ¹		
Delivery Group (SDG)	Sampling Date	Protocol	Sample ID	Matrix	voc	svoc	ALD	PFAS	MET	Noncompliance
	6/24/2021	SW846	IA-3	Air	No	-	-	-	-	VOC: ICV %RSD, LCS %R
2106711	6/24/2021	SW846	AA-3	Air	No	-	-	-	-	VOC: ICV %RSD, LCS %R
	6/24/2021	SW846	DUP-062421	Air	No	-	-	-	-	VOC: ICV %RSD, LCS %R

Note:

Samples which are compliant with no added validation qualifiers are listed as "yes". Samples which are non-compliant or which have added qualifiers are listed as "no". A "no" designation does not necessarily indicate that the data have been rejected or are otherwise unusable.

DATA USABILITY SUMMARY REPORT

VALIDATION PERFORMED BY: Jeffrey L. Davin

SIGNATURE:

DATE: February 28, 2021

PEER REVIEW: Todd Church

DATE: July 30, 2021

CHAIN OF CUSTODY CORRECTED SAMPLE ANALYSIS DATA SHEETS

CULOIUS

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PID: Workorder #: 2106711 Click links below to view: Canister Sampling Guide Helium Shroud Video

Air Toxics against any claim, demand, or action, of any kind, related to the collection, 0) 467-4922	d international law יי, of any kind, rela	rederal, an ind, or action	im, dema	any cla	cs against	rofins Air Toxi ne (800) 467-	t to hold harmless, defend, and indemnify Eurofins Air Toxics aghandling, of shipping of samples. D.O.T Hotline (800) 467-4922	effend, and of samples	armless, de of shipping	ant to hold handling, c	indicates agreeme	ordinances of any kind. Relinquishing signature also indicates agreement to hold harmless, defend, and indemnify Eurofins Air Toxics against any claim, demand, or action, of any kind, related to the collection handling, of shipping of samples. D.O.T Hotline (800) 467-4922	ordinances o
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Client ID: IA-3

Lab ID: 2106711-01A **Date/Time Analyzed:** 7/2/21 03:46 PM

Date/Time Collected: 6/24/21 06:25 PM **Dilution Factor:** 1.42

Media: Instrument/Filename: msd21.i / 21070214

		MDL	LOD	Rpt. Limit	Amount
Compound	CAS#	(ug/m3)	(ug/m3)	(ug/m3)	(ug/m3)
1,2,4-Trichlorobenzene	120-82-1	1.8	2.1	5.3	Not Detected
1,2,4-Trimethylbenzene	95-63-6	0.13	0.28	0.70	Not Detected
1,2-Dichlorobenzene	95-50-1	0.11	0.34	0.85	Not Detected
1,2-Dichloropropane	78-87-5	0.067	0.26	0.66	Not Detected
1,3,5-Trimethylbenzene	108-67-8	0.15	0.28	0.70	Not Detected
1,3-Butadiene	106-99-0	0.049	0.12	0.31	Not Detected
1,3-Dichlorobenzene	541-73-1	0.086	0.34	0.85	Not Detected
1,4-Dioxane	123-91-1	0.074	0.20	0.51	0.14 J
2,2,4-Trimethylpentane	540-84-1	0.21	1.3	3.3	Not Detected
2-Butanone (Methyl Ethyl Ketone)	78-93-3	0.16	0.84	2.1	1.1 J
2-Hexanone	591-78-6	0.16	1.2	2.9	Not Detected
2-Propanol	67-63-0	0.15	0.70	1.7	6.2
3-Chloropropene	107-05-1	0.69	0.89	2.2	Not Detected J
4-Ethyltoluene	622-96-8	0.12	0.28	0.70	Not Detected
4-Methyl-2-pentanone	108-10-1	0.13	0.23	0.58	0.14 J
Acetone	67-64-1	0.55	0.67	3.4	11
alpha-Chlorotoluene	100-44-7	0.12	0.29	0.74	Not Detected
Bromodichloromethane	75-27-4	0.10	0.38	0.95	Not Detected
Bromoform	75-25-2	0.20	0.59	1.5	Not Detected
Bromomethane	74-83-9	0.33	1.1	2.8	Not Detected
Carbon Disulfide	75-15-0	0.39	0.88	2.2	Not Detected
Chlorobenzene	108-90-7	0.061	0.26	0.65	Not Detected
cis-1,3-Dichloropropene	10061-01-5	0.064	0.26	0.64	Not Detected
Cumene	98-82-8	0.074	0.28	0.70	Not Detected



Client ID: IA-3

Lab ID: 2106711-01A **Date/Time Analyzed:** 7/2/21 03:46 PM

Date/Time Collected: 6/24/21 06:25 PM Dilution Factor: 1.42

Media: Instrument/Filename: msd21.i / 21070214

		MDL	LOD	Rpt. Limit	Amount
Compound	CAS#	(ug/m3)	(ug/m3)	(ug/m3)	(ug/m3)
Cyclohexane	110-82-7	0.13	0.98	2.4	Not Detected
Dibromochloromethane	124-48-1	0.15	0.48	1.2	Not Detected
Ethanol	64-17-5	0.14	0.54	1.3	100 J
Freon 11	75-69-4	0.070	0.32	0.80	1.1
Freon 113	76-13-1	0.15	0.44	1.1	0.42 J
Heptane	142-82-5	0.18	1.2	2.9	Not Detected
Hexachlorobutadiene	87-68-3	1.4	3.0	7.6	Not Detected
Hexane	110-54-3	0.15	1.0	2.5	Not Detected
Methylene Chloride	75-09-2	0.80	2.0	0.99	Not Detected
Propylbenzene	103-65-1	0.10	0.28	0.70	Not Detected
Styrene	100-42-5	0.061	0.24	0.60	0.070 J
Tetrahydrofuran	109-99-9	0.21	0.84	2.1	Not Detected
trans-1,3-Dichloropropene	10061-02-6	0.068	0.26	0.64	Not Detected

D: Analyte not within the DoD scope of accreditation.

TENTATIVELY IDENTIFIED COMPOUNDS

	_	Match	LOD	Amount	
Compound	CAS#			ppbv	
1,2,3-Trimethylbenzene	526-73-8	NA		Not Detected	
2-Nitropropane	79-46-9	NA		Not Detected	
4-Chlorotoluene	106-43-4	NA		Not Detected	
Acetonitrile	75-05-8	NA		Not Detected	
Benzaldehyde	100-52-7	NA		Not Detected	
bis(2-Chloroethyl) Ether	111-44-4	NA		Not Detected	



Client ID: IA-3

Lab ID: 2106711-01A **Date/Time Analyzed:** 7/2/21 03:46 PM

Date/Time Collected: 6/24/21 06:25 PM **Dilution Factor:** 1.42

Media: 6 Liter Summa Canister (SIM Certified) Instrument/Filename: msd21.i / 21070214

TENTATIVELY IDENTIFIED COMPOUNDS

		Match	LOD	Amount	
Compound	CAS#			ppbv	
Cyclohexene	110-83-8	NA		Not Detected	
Epichlorohydrin	106-89-8	NA		Not Detected	
Ethyl Acetate	141-78-6	NA		Not Detected	
Isobutanol	78-83-1	NA		Not Detected	
Isopropyl ether	108-20-3	NA		Not Detected	
Methacrylonitrile	126-98-7	NA		Not Detected	
Methyl Acetate	79-20-9	NA		Not Detected	
N,N-Dimethyl Aniline	121-69-7	NA		Not Detected	
n-Butanol	71-36-3	NA		Not Detected	
Nitrobenzene	98-95-3	NA		Not Detected	
Pentane	109-66-0	78%		1.6 NJ	
Propylene	115-07-1	NA		Not Detected	
Pyridine	110-86-1	NA		Not Detected	

J = Estimated value.

D: Analyte not within the DoD scope of accreditation.

Surrogates	CAS#	Limits	%Recovery
1,2-Dichloroethane-d4	17060-07-0	70-130	102
4-Bromofluorobenzene	460-00-4	70-130	85
Toluene-d8	2037-26-5	70-130	96

NJ =The identification is based on presumptive evidence; estimated value.



Client ID: IA-3

Lab ID: 2106711-01B **Date/Time Analyzed:** 7/2/21 03:46 PM

Date/Time Collected: 6/24/21 06:25 PM Dilution Factor: 1.42

Media: 6 Liter Summa Canister (SIM Certified) Instrument/Filename: msd21.i / 21070214sim

		MDL	LOD	Rpt. Limit	Amount
Compound	CAS#	(ug/m3)	(ug/m3)	(ug/m3)	(ug/m3)
1,1,1-Trichloroethane	71-55-6	0.0043	0.062	0.15	0.013 J
1,1,2,2-Tetrachloroethane	79-34-5	0.025	0.078	0.19	Not Detected
1,1,2-Trichloroethane	79-00-5	0.014	0.062	0.15	Not Detected
1,1-Dichloroethane	75-34-3	0.031	0.046	0.11	Not Detected
1,1-Dichloroethene	75-35-4	0.0041	0.045	0.056	Not Detected
1,2-Dibromoethane (EDB)	106-93-4	0.014	0.087	0.22	Not Detected
1,2-Dichloroethane	107-06-2	0.0060	0.046	0.11	0.042 J
1,4-Dichlorobenzene	106-46-7	0.056	0.068	0.17	Not Detected
Benzene	71-43-2	0.013	0.036	0.23	0.11 J
Carbon Tetrachloride	56-23-5	0.0089	0.071	0.18	0.32
Chloroethane	75-00-3	0.0060	0.030	0.19	0.041 J
Chloroform	67-66-3	0.011	0.055	0.14	0.14
Chloromethane	74-87-3	0.18	0.59	1.5	0.74 J
cis-1,2-Dichloroethene	156-59-2	0.0068	0.045	0.11	0.021 J
Ethyl Benzene	100-41-4	0.0046	0.049	0.12	0.058 J
Freon 114	76-14-2	0.017	0.079	0.20	0.092 J
Freon 12	75-71-8	0.0052	0.056	3.5	2.2 J
m,p-Xylene	108-38-3	0.0075	0.049	0.25	0.20 J
Methyl tert-butyl ether	1634-04-4	0.0092	0.041	0.51	Not Detected
Naphthalene	91-20-3	0.094	0.15	0.37	0.43
o-Xylene	95-47-6	0.0086	0.049	0.12	0.071 J
Tetrachloroethene	127-18-4	0.010	0.077	0.19	0.082 J
Toluene	108-88-3	0.0098	0.043	0.27	0.45
trans-1,2-Dichloroethene	156-60-5	0.0058	0.045	0.56	Not Detected



Client ID: IA-3

Lab ID: 2106711-01B **Date/Time Analyzed:** 7/2/21 03:46 PM

Media: 6 Liter Summa Canister (SIM Certified) Instrument/Filename: msd21.i / 21070214sim

Compound	CAS#	MDL (ug/m3)	LOD (ug/m3)	Rpt. Limit (ug/m3)	Amount (ug/m3)
Trichloroethene	79-01-6	0.014	0.061	0.15	0.16
Vinyl Chloride	75-01-4	0.0046	0.029	0.036	Not Detected

J = Estimated value.

D: Analyte not within the DoD scope of accreditation.

Surrogates	CAS#	Limits	%Recovery
1,2-Dichloroethane-d4	17060-07-0	70-130	95
4-Bromofluorobenzene	460-00-4	70-130	88
Toluene-d8	2037-26-5	70-130	94



Client ID: AA-3

Lab ID: 2106711-02A **Date/Time Analyzed:** 7/2/21 04:24 PM

Date/Time Collected: 6/24/21 06:30 PM Dilution Factor: 1.51

Media: Instrument/Filename: msd21.i / 21070215

		MDL	LOD	Rpt. Limit	Amount
Compound	CAS#	(ug/m3)	(ug/m3)	(ug/m3)	(ug/m3)
1,2,4-Trichlorobenzene	120-82-1	1.9	2.2	5.6	Not Detected
1,2,4-Trimethylbenzene	95-63-6	0.14	0.30	0.74	Not Detected
1,2-Dichlorobenzene	95-50-1	0.12	0.36	0.91	Not Detected
1,2-Dichloropropane	78-87-5	0.071	0.28	0.70	Not Detected
1,3,5-Trimethylbenzene	108-67-8	0.16	0.30	0.74	Not Detected
1,3-Butadiene	106-99-0	0.052	0.13	0.33	Not Detected
1,3-Dichlorobenzene	541-73-1	0.092	0.36	0.91	Not Detected
1,4-Dioxane	123-91-1	0.079	0.22	0.54	0.098 J
2,2,4-Trimethylpentane	540-84-1	0.22	1.4	3.5	Not Detected
2-Butanone (Methyl Ethyl Ketone)	78-93-3	0.17	0.89	2.2	1.0 J
2-Hexanone	591-78-6	0.17	1.2	3.1	Not Detected
2-Propanol	67-63-0	0.16	0.74	1.8	2.6
3-Chloropropene	107-05-1	0.74	0.94	2.4	Not Detected J
4-Ethyltoluene	622-96-8	0.12	0.30	0.74	Not Detected
4-Methyl-2-pentanone	108-10-1	0.14	0.25	0.62	Not Detected
Acetone	67-64-1	0.58	0.72	3.6	9.9
alpha-Chlorotoluene	100-44-7	0.13	0.31	0.78	Not Detected
Bromodichloromethane	75-27-4	0.11	0.40	1.0	Not Detected
Bromoform	75-25-2	0.22	0.62	1.6	Not Detected
Bromomethane	74-83-9	0.35	1.2	2.9	Not Detected
Carbon Disulfide	75-15-0	0.42	0.94	2.4	Not Detected
Chlorobenzene	108-90-7	0.065	0.28	0.70	Not Detected
cis-1,3-Dichloropropene	10061-01-5	0.068	0.27	0.68	Not Detected
Cumene	98-82-8	0.078	0.30	0.74	Not Detected



Client ID: AA-3

Lab ID: 2106711-02A **Date/Time Analyzed:** 7/2/21 04:24 PM

Media: Instrument/Filename: msd21.i / 21070215

		MDL	LOD	Rpt. Limit	Amount
Compound	CAS#	(ug/m3)	(ug/m3)	(ug/m3)	(ug/m3)
Cyclohexane	110-82-7	0.13	1.0	2.6	0.21 J
Dibromochloromethane	124-48-1	0.16	0.51	1.3	Not Detected
Ethanol	64-17-5	0.15	0.57	1.4	1.9 J
Freon 11	75-69-4	0.074	0.34	0.85	1.2
Freon 113	76-13-1	0.16	0.46	1.2	0.46 J
Heptane	142-82-5	0.19	1.2	3.1	Not Detected
Hexachlorobutadiene	87-68-3	1.5	3.2	8.0	Not Detected
Hexane	110-54-3	0.16	1.1	2.7	Not Detected
Methylene Chloride	75-09-2	0.85	2.1	1.0	Not Detected
Propylbenzene	103-65-1	0.11	0.30	0.74	Not Detected
Styrene	100-42-5	0.065	0.26	0.64	Not Detected
Tetrahydrofuran	109-99-9	0.22	0.89	2.2	Not Detected
trans-1,3-Dichloropropene	10061-02-6	0.072	0.27	0.68	Not Detected

D: Analyte not within the DoD scope of accreditation.

TENTATIVELY IDENTIFIED COMPOUNDS

	_	Match	LOD	Amount	
Compound	CAS#			ppbv	
1,2,3-Trimethylbenzene	526-73-8	NA		Not Detected	
2-Nitropropane	79-46-9	NA		Not Detected	
4-Chlorotoluene	106-43-4	NA		Not Detected	
Acetonitrile	75-05-8	NA		Not Detected	
Benzaldehyde	100-52-7	NA		Not Detected	
bis(2-Chloroethyl) Ether	111-44-4	NA		Not Detected	



Client ID: AA-3

Lab ID: 2106711-02A **Date/Time Analyzed:** 7/2/21 04:24 PM

Date/Time Collected: 6/24/21 06:30 PM Dilution Factor: 1.51

Media: 6 Liter Summa Canister (SIM Certified) Instrument/Filename: msd21.i / 21070215

TENTATIVELY IDENTIFIED COMPOUNDS

		Match	LOD	Amount	
Compound	CAS#			ppbv	
Cyclohexene	110-83-8	NA		Not Detected	
Epichlorohydrin	106-89-8	NA		Not Detected	
Ethyl Acetate	141-78-6	NA		Not Detected	
Isobutanol	78-83-1	NA		Not Detected	
Isopropyl ether	108-20-3	NA		Not Detected	
Methacrylonitrile	126-98-7	NA		Not Detected	
Methyl Acetate	79-20-9	NA		Not Detected	
N,N-Dimethyl Aniline	121-69-7	NA		Not Detected	
n-Butanol	71-36-3	NA		Not Detected	
Nitrobenzene	98-95-3	NA		Not Detected	
Pentane	109-66-0	NA		Not Detected	
Propylene	115-07-1	NA		Not Detected	
Pyridine	110-86-1	NA		Not Detected	

J = Estimated value.

D: Analyte not within the DoD scope of accreditation.

Surrogates	CAS#	Limits	%Recovery
1,2-Dichloroethane-d4	17060-07-0	70-130	99
4-Bromofluorobenzene	460-00-4	70-130	83
Toluene-d8	2037-26-5	70-130	99



Client ID: AA-3

Lab ID: 2106711-02B **Date/Time Analyzed:** 7/2/21 04:24 PM

Date/Time Collected: 6/24/21 06:30 PM Dilution Factor: 1.51

Media: 6 Liter Summa Canister (SIM Certified) Instrument/Filename: msd21.i / 21070215sim

		MDL	LOD	Rpt. Limit	Amount
Compound	CAS#	(ug/m3)	(ug/m3)	(ug/m3)	(ug/m3)
1,1,1-Trichloroethane	71-55-6	0.0045	0.066	0.16	0.0082 J
1,1,2,2-Tetrachloroethane	79-34-5	0.026	0.083	0.21	Not Detected
1,1,2-Trichloroethane	79-00-5	0.015	0.066	0.16	Not Detected
1,1-Dichloroethane	75-34-3	0.033	0.049	0.12	Not Detected
1,1-Dichloroethene	75-35-4	0.0044	0.048	0.060	Not Detected
1,2-Dibromoethane (EDB)	106-93-4	0.015	0.093	0.23	Not Detected
1,2-Dichloroethane	107-06-2	0.0064	0.049	0.12	0.049 J
1,4-Dichlorobenzene	106-46-7	0.060	0.073	0.18	Not Detected
Benzene	71-43-2	0.014	0.038	0.24	0.10 J
Carbon Tetrachloride	56-23-5	0.0095	0.076	0.19	0.40
Chloroethane	75-00-3	0.0064	0.032	0.20	0.030 J
Chloroform	67-66-3	0.012	0.059	0.15	0.065 J
Chloromethane	74-87-3	0.19	0.62	1.6	0.71 J
cis-1,2-Dichloroethene	156-59-2	0.0072	0.048	0.12	Not Detected
Ethyl Benzene	100-41-4	0.0048	0.052	0.13	0.044 J
Freon 114	76-14-2	0.018	0.084	0.21	0.10 J
Freon 12	75-71-8	0.0055	0.060	3.7	2.1 J
m,p-Xylene	108-38-3	0.0080	0.052	0.26	0.15 J
Methyl tert-butyl ether	1634-04-4	0.0098	0.044	0.54	Not Detected
Naphthalene	91-20-3	0.10	0.16	0.40	0.12 J
o-Xylene	95-47-6	0.0092	0.052	0.13	0.052 J
Tetrachloroethene	127-18-4	0.011	0.082	0.20	0.048 J
Toluene	108-88-3	0.010	0.046	0.28	0.24 J
trans-1,2-Dichloroethene	156-60-5	0.0062	0.048	0.60	Not Detected



Client ID: AA-3

Lab ID: 2106711-02B **Date/Time Analyzed:** 7/2/21 04:24 PM

Date/Time Collected: 6/24/21 06:30 PM Dilution Factor: 1.51

Media: 6 Liter Summa Canister (SIM Certified) Instrument/Filename: msd21.i / 21070215sim

Compound	CAS#	MDL (ug/m3)	LOD (ug/m3)	Rpt. Limit (ug/m3)	Amount (ug/m3)
Trichloroethene	79-01-6	0.015	0.065	0.16	0.017 J
Vinyl Chloride	75-01-4	0.0049	0.031	0.038	Not Detected

J = Estimated value.

D: Analyte not within the DoD scope of accreditation.

Surrogates	CAS#	Limits	%Recovery
1,2-Dichloroethane-d4	17060-07-0	70-130	94
4-Bromofluorobenzene	460-00-4	70-130	84
Toluene-d8	2037-26-5	70-130	93



 Client ID:
 DUP-062421

 Lab ID:
 2106711-03A

Date/Time Collected: 6/24/21 12:00 AM

Media:

Date/Time Analyzed: 7/2/21 05:01 PM

Dilution Factor: 1.39

Instrument/Filename: msd21.i / 21070216

		MDL	LOD	Rpt. Limit	Amount
Compound	CAS#	(ug/m3)	(ug/m3)	(ug/m3)	(ug/m3)
1,2,4-Trichlorobenzene	120-82-1	1.7	2.1	5.2	Not Detected
1,2,4-Trimethylbenzene	95-63-6	0.12	0.27	0.68	Not Detected
1,2-Dichlorobenzene	95-50-1	0.11	0.33	0.84	Not Detected
1,2-Dichloropropane	78-87-5	0.065	0.26	0.64	Not Detected
1,3,5-Trimethylbenzene	108-67-8	0.15	0.27	0.68	Not Detected
1,3-Butadiene	106-99-0	0.048	0.12	0.31	Not Detected
1,3-Dichlorobenzene	541-73-1	0.084	0.33	0.84	Not Detected
1,4-Dioxane	123-91-1	0.072	0.20	0.50	0.19 J
2,2,4-Trimethylpentane	540-84-1	0.20	1.3	3.2	Not Detected
2-Butanone (Methyl Ethyl Ketone)	78-93-3	0.15	0.82	2.0	1.4 J
2-Hexanone	591-78-6	0.16	1.1	2.8	0.19 J
2-Propanol	67-63-0	0.15	0.68	1.7	6.4
3-Chloropropene	107-05-1	0.68	0.87	2.2	Not Detected J
4-Ethyltoluene	622-96-8	0.11	0.27	0.68	Not Detected
4-Methyl-2-pentanone	108-10-1	0.13	0.23	0.57	0.15 J
Acetone	67-64-1	0.54	0.66	3.3	12
alpha-Chlorotoluene	100-44-7	0.12	0.29	0.72	Not Detected
Bromodichloromethane	75-27-4	0.10	0.37	0.93	Not Detected
Bromoform	75-25-2	0.20	0.57	1.4	Not Detected
Bromomethane	74-83-9	0.32	1.1	2.7	Not Detected
Carbon Disulfide	75-15-0	0.38	0.86	2.2	0.65 J
Chlorobenzene	108-90-7	0.060	0.26	0.64	Not Detected
cis-1,3-Dichloropropene	10061-01-5	0.062	0.25	0.63	Not Detected
Cumene	98-82-8	0.072	0.27	0.68	Not Detected



Client ID: DUP-062421 **Lab ID:** 2106711-03A

Date/Time Collected: 6/24/21 12:00 AM

Media:

Date/Time Analyzed: 7/2/21 05:01 PM

Dilution Factor: 1.39

Instrument/Filename: msd21.i / 21070216

		MDL	LOD	Rpt. Limit	Amount
Compound	CAS#	(ug/m3)	(ug/m3)	(ug/m3)	(ug/m3)
Cyclohexane	110-82-7	0.12	0.96	2.4	1.6 J
Dibromochloromethane	124-48-1	0.14	0.47	1.2	Not Detected
Ethanol	64-17-5	0.14	0.52	1.3	100 ^J
Freon 11	75-69-4	0.068	0.31	0.78	1.1
Freon 113	76-13-1	0.15	0.43	1.1	0.38 J
Heptane	142-82-5	0.18	1.1	2.8	Not Detected
Hexachlorobutadiene	87-68-3	1.4	3.0	7.4	Not Detected
Hexane	110-54-3	0.14	0.98	2.4	Not Detected
Methylene Chloride	75-09-2	0.78	1.9	0.96	Not Detected
Propylbenzene	103-65-1	0.10	0.27	0.68	Not Detected
Styrene	100-42-5	0.060	0.24	0.59	0.093 J
Tetrahydrofuran	109-99-9	0.20	0.82	2.0	0.22 J
trans-1,3-Dichloropropene	10061-02-6	0.066	0.25	0.63	Not Detected

D: Analyte not within the DoD scope of accreditation.

TENTATIVELY IDENTIFIED COMPOUNDS

		Match	LOD	Amount	
Compound	CAS#			ppbv	
1,2,3-Trimethylbenzene	526-73-8	NA		Not Detected	
2-Nitropropane	79-46-9	NA		Not Detected	
4-Chlorotoluene	106-43-4	NA		Not Detected	
Acetonitrile	75-05-8	NA		Not Detected	
Benzaldehyde	100-52-7	NA		Not Detected	
bis(2-Chloroethyl) Ether	111-44-4	NA		Not Detected	



Client ID: DUP-062421

Lab ID: 2106711-03A Date/Time Analyzed:

Date/Time Collected: 6/24/21 12:00 AM **Dilution Factor:** 1.39

Media: 6 Liter Summa Canister (SIM Certified) Instrument/Filename: msd21.i / 21070216

TENTATIVELY IDENTIFIED COMPOUNDS

7/2/21 05:01 PM

		Match	LOD	Amount	
Compound	CAS#			ppbv	
Cyclohexene	110-83-8	NA		Not Detected	
Epichlorohydrin	106-89-8	NA		Not Detected	
Ethyl Acetate	141-78-6	NA		Not Detected	
Isobutanol	78-83-1	NA		Not Detected	
Isopropyl ether	108-20-3	NA		Not Detected	
Methacrylonitrile	126-98-7	NA		Not Detected	
Methyl Acetate	79-20-9	NA		Not Detected	
N,N-Dimethyl Aniline	121-69-7	NA		Not Detected	
n-Butanol	71-36-3	NA		Not Detected	
Nitrobenzene	98-95-3	NA		Not Detected	
Pentane	109-66-0	86%		1.6 NJ	
Propylene	115-07-1	NA		Not Detected	
Pyridine	110-86-1	NA		Not Detected	

J = Estimated value.

D: Analyte not within the DoD scope of accreditation.

Surrogates	CAS#	Limits	%Recovery
1,2-Dichloroethane-d4	17060-07-0	70-130	98
4-Bromofluorobenzene	460-00-4	70-130	80
Toluene-d8	2037-26-5	70-130	94

NJ =The identification is based on presumptive evidence; estimated value.



Client ID: DUP-062421 **Lab ID:** 2106711-03B

Date/Time Collected: 6/24/21 12:00 AM

Media: 6 Liter Summa Canister (SIM Certified)

Date/Time Analyzed: 7/2/21 05:01 PM

Dilution Factor: 1.39

Instrument/Filename: msd21.i / 21070216sim

		MDL	LOD	Rpt. Limit	Amount
Compound	CAS#	(ug/m3)	(ug/m3)	(ug/m3)	(ug/m3)
1,1,1-Trichloroethane	71-55-6	0.0042	0.061	0.15	0.015 J
1,1,2,2-Tetrachloroethane	79-34-5	0.024	0.076	0.19	Not Detected
1,1,2-Trichloroethane	79-00-5	0.014	0.061	0.15	Not Detected
1,1-Dichloroethane	75-34-3	0.031	0.045	0.11	Not Detected
1,1-Dichloroethene	75-35-4	0.0040	0.044	0.055	Not Detected
1,2-Dibromoethane (EDB)	106-93-4	0.014	0.085	0.21	Not Detected
1,2-Dichloroethane	107-06-2	0.0058	0.045	0.11	0.043 J
1,4-Dichlorobenzene	106-46-7	0.055	0.067	0.17	Not Detected
Benzene	71-43-2	0.012	0.036	0.22	0.12 J
Carbon Tetrachloride	56-23-5	0.0087	0.070	0.17	0.33
Chloroethane	75-00-3	0.0059	0.029	0.18	0.026 J
Chloroform	67-66-3	0.011	0.054	0.14	0.15
Chloromethane	74-87-3	0.17	0.57	1.4	0.74 J
cis-1,2-Dichloroethene	156-59-2	0.0067	0.044	0.11	0.022 J
Ethyl Benzene	100-41-4	0.0045	0.048	0.12	0.065 J
Freon 114	76-14-2	0.017	0.078	0.19	0.095 J
Freon 12	75-71-8	0.0051	0.055	3.4	2.2 J
m,p-Xylene	108-38-3	0.0074	0.048	0.24	0.22 J
Methyl tert-butyl ether	1634-04-4	0.0090	0.040	0.50	Not Detected
Naphthalene	91-20-3	0.092	0.14	0.36	0.42
o-Xylene	95-47-6	0.0084	0.048	0.12	0.093 J
Tetrachloroethene	127-18-4	0.0098	0.075	0.19	0.060 J
Toluene	108-88-3	0.0096	0.042	0.26	0.51
trans-1,2-Dichloroethene	156-60-5	0.0057	0.044	0.55	Not Detected



Client ID: DUP-062421 **Lab ID:** 2106711-03B

Date/Time Collected: 6/24/21 12:00 AM

Media: 6 Liter Summa Canister (SIM Certified)

Date/Time Analyzed: 7/2/21 05:01 PM

Dilution Factor: 1.39

Instrument/Filename: msd21.i / 21070216sim

		MDL	LOD	Rpt. Limit	Amount
Compound	CAS#	(ug/m3)	(ug/m3)	(ug/m3)	(ug/m3)
Trichloroethene	79-01-6	0.014	0.060	0.15	0.15
Vinyl Chloride	75-01-4	0.0045	0.028	0.036	Not Detected

J = Estimated value.

D: Analyte not within the DoD scope of accreditation.

Surrogates	CAS#	Limits	%Recovery
1,2-Dichloroethane-d4	17060-07-0	70-130	95
4-Bromofluorobenzene	460-00-4	70-130	85
Toluene-d8	2037-26-5	70-130	94



Bristol Myers Squibb Thompson Road Investigation

Data Usability Summary Report

Syracuse, NY

Volatile Organic Compound (VOC) Analysis

SDGs # 2104655, 2105127, 2105279A and 2105279B

Analyses Performed By: Eurofins Air Toxics

Folsom, CA

Report #41885R Review Level: Tier III Project: 30064943

SUMMARY

This data quality assessment summarizes the review of Sample Delivery Groups (SDGs) # 2104655, 2105127, 2105279A and 2105279B for samples collected in association with the Bristol Myers Squibb Thompson Road Site. The review was conducted as a Tier III evaluation and included review of data package completeness. Only analytical data associated with constituents of concern were reviewed for this validation. Field documentation was not included in this review. Included with this assessment are the validation annotated sample result sheets, and chain of custody. Analyses were performed on the following samples:

SDG				Sample	Parent		Analysis			
Number	Sample ID	Lab ID	Matrix	Collection Date	Sample	voc	svoc	PFAS	MET	ALD
2104655	IA-3_042921	2104655-01	Air	4/29/2021		Х				
2105127	IA-3_050621	2105127-01	Air	5/6/2021		Х				
2105279A	IA-3_051321	2105279A-01	Air	5/13/2021		Х				
2105279B	IA-3_051321	2105279B-01	Air	5/13/2021		Х				

Note: SDG 2105279B is a reanalysis of the sample in 2105279A to meet the required action limits for Trichloroethene.

ANALYTICAL DATA PACKAGE DOCUMENTATION

The table below is the evaluation of the data package completeness.

	Rep	Reported		rmance ptable	Not
Items Reviewed	No	Yes	No	Yes	Required
Sample receipt condition		Х		Х	
2. Requested analyses and sample results		Х		Х	
Master tracking list		Х		Х	
4. Methods of analysis		Х		Х	
5. Reporting limits		Х		Х	
6. Sample collection date		Х		Х	
7. Laboratory sample received date		Х		Х	
8. Sample preservation verification (as applicable)		Х		Х	
9. Sample preparation/extraction/analysis dates		Х		Х	
10. Fully executed Chain-of-Custody (COC) form		Х		Х	
11. Narrative summary of QA or sample problems provided		Х		X	
12. Data Package Completeness and Compliance		Х		Х	

Note:

QA - Quality Assurance

List of Acronyms

%D: Percent Difference

%R: Percent Recovery

AC: Acceptable

ALC/GLY: Alcohols/Glycols

BAL: Blank Action Level

CCV: Continuing Calibration Verification

CRDL: Contract Required Detection Limit

D: Dilution

EIS: Extractable Internal Standard

FB: Field Blank

FD: Field Duplicate

ALD: Aldehydes

GC/ECD: Gas Chromatograph/Electron Capture Detector

GC/MS: Gas Chromatograph/Mass spectrometer

HT: Holding Time

ICP: Inductively Coupled Plasma

ICS: Interference Control Sample

ICV: Initial Calibration Verification

ISTD: Internal Standards

LabDup: Laboratory Duplicate

LCS: Lab Control Sample

LCSD: Lab Control Sample Duplicate

LL: Lower Control Limit

MB: Method Blank

MDL: Method Detection Limit

MET: Metals

MS: Matrix Spike

MSD: Matrix Spike Duplicate

N/A: Not Applicable

NC: Not Compliant

List of Acronyms, Continued

PAH: Polyaromatic Hydrocarbon

PCB: Polychlorinated Biphenyl

PEST: Pesticide

PFAS: Per- and Polyfluoroalkyl Substances

QA: Quality Assurance

QC: Quality Control

RB: Rinse Blank

RL: Reporting Limit

RPD: Relative Percent Difference

RRF: Relative Response Factor

RSD: Relative Standard Deviation

RT: Retention Time

SDG: Sample Delivery Group

SerDil: Serial Dilution

SIM: Single Ion Monitoring

SOP: Standard Operating Procedure

SSTD: Surrogate Standards

SVOC: Semivolatile Organic Compound

TB: Trip Blank

TIC: Tentatively Identified Compound

TOC: Total Organic Carbon

TOTDIS: Total and Dissolved

UL: Upper Control Limit

USEPA: United States Environmental Protection Agency

VOC: Volatile Organic Compound

ORGANIC ANALYSIS INTRODUCTION

Analyses were performed according to United States Environmental Protection Agency (USEPA) Method TO-15. Data were reviewed in accordance with USEPA National Functional Guidelines of October 1999, USEPA Region II SOP HW-31- Validating Air Samples Volatile Organic Analysis of Ambient Air In Canister by Method TO-15 of October 2006, New York State DEC Analytical Method ASP 2005 TO-15 (QA/QC Criteria R9 TO-15), and NYSDEC Modifications to R9 TO-15 QA/QC Criteria October 2009.

The data review process is an evaluation of data on a technical basis rather than a determination of contract compliance. As such, the standards against which the data are being weighed may differ from those specified in the analytical method. It is assumed that the data package represents the best efforts of the laboratory and had already been subjected to adequate and sufficient quality review prior to submission.

During the review process, laboratory qualified and unqualified data are verified against the supporting documentation. Based on this evaluation, qualifier codes may be added, deleted, or modified by the data reviewer. Results are qualified with the following codes in accordance with USEPA National Functional Guidelines:

- Concentration (C) Qualifiers
 - U The compound was analyzed for but not detected. The associated value is the compound quantitation limit.
 - B The compound has been found in the sample as well as its associated blank, its presence in the sample may be suspect.
- Quantitation (Q) Qualifiers
 - E The compound was quantitated above the calibration range.
 - D Concentration is based on a diluted sample analysis.
- Validation Qualifiers
 - J The analyte was positively identified; however, the associated numerical value is an estimated concentration only.
 - UJ The analyte was not detected above the reported sample detection limit. However, the reported limit is approximate and may or may not represent the actual limit of detection.
 - UB Analyte considered non-detect at the listed value due to associated blank contamination.
 - N The analysis indicates the presence of a compound for which there is presumptive evidence to make a tentative identification.
 - R The sample results are rejected.

Two facts should be noted by all data users. First, the "R" flag means that the associated value is unusable. In other words, due to significant quality control (QC) problems, the analysis is invalid and provides no information as to whether the compound is present or not. "R" values should not appear on data tables because they cannot be relied upon, even as a last resort. The second fact to keep in mind is that no compound concentration, even if it has passed all QC tests, is guaranteed to be accurate. Strict QC serves to increase confidence in data but any value potentially contains error.

VOLATILE ORGANIC COMPOUND (VOC) ANALYSES

1. Holding Times

The specified holding times for the following methods are presented in the following table.

Method	Matrix	Holding Time	Preservation	Return Canister Pressure
USEPA TO-15	Air	30 days from collection to analysis	Ambient Temperature	< -1" Hg

All samples were analyzed within the specified holding time and canister return pressure / vacuum criteria.

2. Blank Contamination

Quality assurance (QA) blanks (i.e., method and rinse blanks) are prepared to identify any contamination which may have been introduced into the samples during sample preparation or field activity. Method blanks measure laboratory contamination. Rinse blanks measure contamination of samples during field operations.

A blank action level (BAL) of five times the concentration of a detected compound in an associated blank (common laboratory contaminant compounds are calculated at ten times) is calculated for QA blanks containing concentrations greater than the method detection limit (MDL). The BAL is compared to the associated sample results to determine the appropriate qualification of the sample results, if needed.

All compounds associated with the QA blanks exhibited a concentration less than the MDL, with the exception of the compounds listed in the following table. Sample results associated with QA blank contamination that were greater than the BAL resulted in the removal of the laboratory qualifier (B) of data. Sample results less than the BAL associated with the following sample locations were qualified as listed in the following table.

Sample Locations	Analytes	Sample Result	Qualification
IA-3_051321	Trichloroethene	Detected sample results <rl <bal<="" and="" td=""><td>"UB" at the RL</td></rl>	"UB" at the RL

Note:

RL Reporting limit

3. Mass Spectrometer Tuning

Mass spectrometer performance was acceptable and all analyses were performed within a 12-hour tune clock.

System performance and column resolution were acceptable.

4. Calibration

Satisfactory instrument calibration is established to ensure that the instrument is capable of producing acceptable quantitative data. An initial calibration demonstrates that the instrument is capable of acceptable performance at the beginning of an experimental sequence. The continuing calibration verifies that the instrument daily performance is satisfactory.

4.1 Initial Calibration

The method specifies percent relative standard deviation (%RSD) and relative response factor (RRF) limits for select compounds only. A technical review of the data applies limits to all compounds with no exceptions.

All target compounds associated with the initial calibration standards must exhibit a %RSD less than the control limit (30%) or a correlation coefficient greater than 0.99 and an RRF value greater than control limit (0.05).

4.2 Continuing Calibration

All target compounds associated with the continuing calibration standard must exhibit a percent difference (%D) less than the control limit (30%) and RRF value greater than control limit (0.05).

Compounds associated with the calibrations were within the specified control limits, with the exception of the compounds presented in the following table.

Sample Locations	Initial/Continuing	Compound	Criteria
IA-3_042921	ICV %RSD	Bromomethane	32.9%
IA-3_051321			
	ICV %RSD Freon 11		33.1%
IA-3_050621	CCV %D	Bromomethane	-33.1%
	CCV %D	Carbon Tetrachloride	39.0%
IA-3_051321	CCV %D	Carbon Tetrachloride	30.7%

The criteria used to evaluate the initial and continuing calibration are presented in the following table. In the case of a calibration deviation, the sample results are qualified.

Initial/Continuing	Criteria	Sample Result	Qualification
	RRF <0.05	Non-detect	R
Initial and Continuing Calibration	KKF 40.03	Detect	J
	RRF <0.01 ¹	Non-detect	R
	KKF <0.01	Detect	J
	RRF >0.05 or RRF >0.01 ¹	Non-detect	No Action

Initial/Continuing	Criteria	Sample Result	Qualification
		Detect	
	%RSD > 15% or a correlation coefficient <0.99	Non-detect	UJ
Initial Calibration	70KSD > 15 % of a correlation coefficient <0.99	coefficient <0.99 Detect J	
	%RSD >90%	Non-detect	R
	70K3D >90%	Detect	J
	9/D > 209/ (increase in consistivity)	Non-detect	No Action
	%D >20% (increase in sensitivity)	Detect	J
Continuing Colibration	9/D > 209/ (degrees in consitiutiv)	Non-detect	UJ
Continuing Calibration	%D >20% (decrease in sensitivity)	Detect	J
	9/ D > 909/ (increase/degrages in consitivity)	Non-detect	R
	%D >90% (increase/decrease in sensitivity)	Detect	J

Note:

5. Surrogates/System Monitoring Compounds

All samples to be analyzed for organic compounds are spiked with surrogate compounds prior to sample preparation to evaluate overall laboratory performance and efficiency of the analytical technique. VOC analysis requires that all surrogates associated with the analysis exhibit a percent recovery within the established acceptance limits of 70% to 130%.

Surrogate recoveries were within control limits.

6. Internal Standard Performance

Internal standard performance criteria ensure that the GC/MS sensitivity and response are stable during every sample analysis. The criteria requires the internal standard compounds associated with the VOC exhibit area counts that are not greater than two times (+100%) or less than one-half (-50%) of the area counts of the associated continuing calibration standard.

Internal standard responses were within control limits.

7. Laboratory Control Sample (LCS) Analysis

The LCS analysis is used to assess the accuracy of the analytical method independent of matrix interferences. The compounds associated with the LCS/LCSD analysis must exhibit a percent recovery within the established acceptance limits of 70% to 130%.

Sample locations associated with LCS/LCSD analysis exhibiting recoveries outside of the control limits presented in the following table.

¹ RRF of 0.01 only applies to compounds which are typically poor responding compounds (i.e., ketones, 1,4-dioxane, etc.)

Sample Locations	Compound	LCS Recovery	LCSD Recovery
14.0.05004	Carbon Tetrachloride	<ll but="">10%</ll>	<ll but="">10%</ll>
IA-3_050621	Hexachlorobutadiene		AC

The criteria used to evaluate the LCS/LCSD recoveries are presented in the following table. In the case of an LCS/LCSD deviation, the sample results are qualified as documented in the table below.

Control Limit	Sample Result	Qualification
the upper central limit (LIL)	Non-detect	No Action
> the upper control limit (UL)	Detect	J
< the lower control limit (LL) but > 10%	Non-detect	UJ
the lower control limit (LL) but > 10%	Detect	J
< 10%	Non-detect	R
1070	Non-detect No Action Detect J Non-detect UJ Detect J	J

8. Field Duplicate Analysis

Field duplicate analysis is used to assess the overall precision of the field sampling procedures and analytical method. A control limit of 50% for air matrices is applied to the RPD between the parent sample and the field duplicate. In the instance when the parent and/or duplicate sample concentrations are less than or equal to 5 times the RL, a control limit of three times the RL is applied for air matrices.

A field duplicate was not included with this SDG.

9. Compound Identification

Compounds are identified on the GC/MS by using the analytes relative retention time and ion spectra.

Sample results associated with compound that exhibited a concentration greater than the linear range of the instrument calibration are summarized in the following table.

Sample ID	Compound	Original Analysis	Diluted Analysis	Reported Analysis
IA-3_042921	Ethanol	-	3000 E	3000 EJ
IA-3_050621	Ethanol	-	1600 E	1600 EJ
IA-3_051321	Ethanol	-	3300 E	3300 EJ

Note: In the instance where both the original analysis and the diluted analysis sample results exhibited a concentration greater than and/or less than the calibration linear range of the instrument; the sample result exhibiting the greatest concentration will be reported as the final result.

Sample results associated with compounds exhibiting concentrations greater than the linear range are qualified as documented in the table below when reported as the final reported sample result.

Reported Sample Results	Qualification
Diluted sample result within calibration range	D
Diluted sample result less than the calibration range	DJ
Diluted sample result greater than the calibration range	EDJ
Original sample result greater than the calibration range	EJ

10. System Performance and Overall Assessment

Please note, the laboratory includes a Limit of Detection (LOD) in the laboratory report which is specific to Department of Defense (DOD) reporting and should not be considered for this site/project. Only the Reporting Limit (RL) and Method Detection Limit (MDL) are stored in the database for this data set.

SDG 2105279B is a reanalysis of the sample in 2105279A to meet the required action limits for Trichloroethene. The result for Trichloroethene associated with IA-3_051321 is reported from SDG 2105279B.

Overall system performance was acceptable. Other than for those deviations specifically mentioned in this review, the overall data quality is within the guidelines specified in the method.

DATA VALIDATION CHECKLIST FOR VOCs

VOCs: TO-15	Re	ported		ormance eptable	Not
	No	Yes	No	Yes	Required
GAS CHROMATOGRAPHY/MASS SPECTROMET	RY (GC/N	MS)			
Tier II Validation					
Holding times		X		X	
Canister return pressure (<-1"Hg)		Х		Х	
Reporting limits (units)		Х		Х	
Blanks		'			
A. Method blanks		Х	Х		
B. Equipment blanks	Х				Х
C. Trip blanks	Х				Х
Laboratory Control Sample (LCS)		Х	Х		
Laboratory Control Sample Duplicate (LCSD)		Х	Х		
LCS/LCSD Precision (RPD)		Х		X	
Matrix Spike (MS)	Х				Х
Matrix Spike Duplicate (MSD)	Х				Х
MS/MSD Precision (RPD)	Х				Х
Field/Lab Duplicate (RPD)	X				Х
Surrogate Spike Recoveries		Х		X	
Dilution Factor		Х		X	
Moisture Content	Х				Х
Tier III Validation					I
System performance and column resolution		X		X	
Initial calibration %RSDs		Х	Х		
Continuing calibration RRFs		X		X	
Continuing calibration %Ds		X	Х		
Instrument tune and performance check		X		X	
Ion abundance criteria for each instrument used		X		X	
Internal standard		X		X	
Compound identification and quantitation					
A. Reconstructed ion chromatograms		Х		Х	

VOCs: TO-15	Rep	oorted		rmance ptable	Not
	No	Yes	No	Yes	Required
GAS CHROMATOGRAPHY/MASS SPECTROMETR	Y (GC/M	S)			
B. Quantitation Reports		Х		X	
C. RT of sample compounds within the established RT windows		Х		Х	
D. Transcription/calculation errors present		X		X	
E. Reporting limits adjusted to reflect sample dilutions		Х		X	

Notes:

%RSD Relative standard deviation

%R Percent recovery

RPD Relative percent difference

%D Percent difference

DATA USABILITY SUMMARY REPORT

SAMPLE COMPLIANCE REPORT

Sample						(Compliar	ıcy¹		
Delivery Group (SDG)	Sampling Date	Protocol	Sample ID	Matrix	voc	svoc	ALD	PFAS	MET	Noncompliance
2104655	4/29/2021	SW846	IA-3_042921	Air	No	-	-	-	-	VOC: ICV %RSD
2105127	5/6/2021	SW846	IA-3_050621	Air	No	-	-	-	-	VOC: CCV %D
2105279A	5/13/2021	SW846	IA-3_051321	Air	No					VOC: ICV %RSD
2105279B	5/13/2021	SW846	IA-3_051321	Air	No	-	-	-	-	VOC: ICV %RSD, MB

Note:

Samples which are compliant with no added validation qualifiers are listed as "yes". Samples which are non-compliant or which have added qualifiers are listed as "no". A "no" designation does not necessarily indicate that the data have been rejected or are otherwise unusable.

DATA USABILITY SUMMARY REPORT

VALIDATION PERFORMED BY: Jeffrey L. Davin

SIGNATURE:

DATE: June 24, 2021

PEER REVIEW: Dennis K. Capria

DATE: June 29, 2021

CHAIN OF CUSTODY CORRECTED SAMPLE ANALYSIS DATA SHEETS

SE CUROTINS	Alialysis Kequest/Vallister Chain o	Senther	Magnisus.	नाजा ह		Gustouy	Ę				· · · ·
Air Toxics 180 Blue Ravine Rd. Suite B, Folsom, CA 95630	PID:	For Laborato	y Use Onl	2104655	ļ.		o o	lick links	Click links below to view: Canister Sampting Guide	.we	
(800) 985-59							}*****	tetium Sh	Helium Shroud Video	K	
Arge	PID	Special In:	Special Instructions/Notes:	<u>-</u>		, ,	n <u>ı</u>	naround	Time (Rush	surcharg	Turnaround Time (Rush surcharges may apply)
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Sample Transportation Notice Bolleville	Fedex Custody Seals Intact?	t? Yes) No	Along	MESTER						
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		handling, of ship	handling, of shipping of samples. D.O.T Hotline (800) 467-4922	.O.T Hotline	(800) 467-49	922					





Air Toxics

MODIFIED EPA METHOD TO-15 GC/MS SIM/FULL SCAN BMS Bldg 3

1.45 msdv.i / v050309 5/3/21 01:01 PM Instrument/Filename: Date/Time Analyzed: Dilution Factor: Date/Time Collected: 4/29/21 04:10 PM 2104655-01A Client ID: Lab ID: Media:

		MDL	ГОБ	Rpt. Limit	Amount
Compound	CAS#	(ng/m3)	(ng/m3)	(ng/m3)	(ng/m3)
1,2,4-Trichlorobenzene	120-82-1	2.4	4.3	5.4	Not Detected
1,2,4-Trimethylbenzene	92-63-6	0.30	0.57	0.71	Not Detected
1,2-Dichlorobenzene	95-50-1	0.52	0.70	0.87	Not Detected
1,2-Dichloropropane	78-87-5	0.31	0.54	0.67	Not Detected
1,3,5-Trimethylbenzene	108-67-8	0.36	0.57	0.71	Not Detected
1,3-Butadiene	106-99-0	0.13	0.26	0.32	Not Detected
1,3-Dichlorobenzene	541-73-1	0.26	0.70	0.87	Not Detected
1,4-Dioxane	123-91-1	0.29	0.42	0.52	0.36 J
2,2,4-Trimethylpentane	540-84-1	0.75	2.7	3.4	Not Detected
2-Butanone (Methyl Ethyl Ketone)	78-93-3	0.35	1.7	2.1	1.5 J
2-Hexanone	591-78-6	0.43	2.4	3.0	Not Detected
2-Propanol	67-63-0	0.39	1.4	1.8	130
3-Chloropropene	107-05-1	1.1	1.8	2.3	Not Detected
4-Ethyltoluene	622-96-8	0.27	0.57	0.71	Not Detected
4-Methyl-2-pentanone	108-10-1	0.13	0.48	0.59	Not Detected
Acetone	67-64-1	0.62	1.4	3.4	22
alpha-Chlorotoluene	100-44-7	0.46	09:0	0.75	Not Detected
Bromodichloromethane	75-27-4	0:30	0.78	0.97	Not Detected
Bromoform	75-25-2	0.46	1.2	1.5	Not Detected
Bromomethane	74-83-9	0.55	2.2	2.8	Not Detected J
Carbon Disulfide	75-15-0	0.33	1.8	2.2	Not Detected
Chlorobenzene	108-90-7	0.21	0.53	0.67	Not Detected
cis-1,3-Dichloropropene	10061-01-5	0.25	0.53	99.0	Not Detected
Cumene	98-82-8	0.094	0.57	0.71	Not Detected

Page 5 of 26





Air Toxics

MODIFIED EPA METHOD TO-15 GC/MS SIM/FULL SCAN BMS Bldg 3

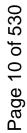
1.45 msdv.i / v050309 5/3/21 01:01 PM Instrument/Filename: Date/Time Analyzed: Dilution Factor: Date/Time Collected: 4/29/21 04:10 PM 2104655-01A Client ID: Lab ID: Media:

		MDL	ГОР	Rpt. Limit	Amount
Compound	CAS#	(ng/m3)	(ng/m3)	(ng/m3)	(ng/m3)
Cyclohexane	110-82-7	0.20	0.40	2.5	Not Detected
Dibromochloromethane	124-48-1	0.49	0.99	1.2	Not Detected
Ethanol	64-17-5	0.39	1.1	1.4	3000 E J
Freon 11	75-69-4	0.18	0.65	0.81	1.3
Freon 113	76-13-1	0.46	0.89	1.1	Not Detected
Heptane	142-82-5	0.55	2.4	3.0	Not Detected
Hexachlorobutadiene	87-68-3	2.4	6.2	7.7	Not Detected
Hexane	110-54-3	0.43	2.0	2.6	Not Detected
Methylene Chloride	75-09-2	0.25	0.40	1.0	0.69 J
Propylbenzene	103-65-1	0.42	0.57	0.71	Not Detected
Styrene	100-42-5	0.18	0.49	0.62	0.52 J
Tetrahydrofuran	109-99-9	0.73	1.7	2.1	0.93 J
trans-1,3-Dichloropropene	10061-02-6	0.28	0.53	99.0	Not Detected

D: Analyte not within the DoD scope of accreditation.

TENTATIVELY IDENTIFIED COMPOUNDS

		Match	ГОР	Amount
Compound	CAS#			ppbv
1,2,3-Trimethylbenzene	526-73-8	AN		Not Detected
2-Nitropropane	79-46-9	Ϋ́		Not Detected
4-Chlorotoluene	106-43-4	Ϋ́		Not Detected
Acetonitrile	75-05-8	Ϋ́		Not Detected
Benzaldehyde	100-52-7	Ϋ́		Not Detected
bis(2-Chloroethyl) Ether	111-44-4	Ϋ́		Not Detected





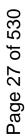
1.45 msdv.i / v050309 5/3/21 01:01 PM Instrument/Filename: Date/Time Analyzed: Dilution Factor: 6 Liter Summa Canister (SIM Certified) Date/Time Collected: 4/29/21 04:10 PM 2104655-01A Client ID: Lab ID: Media:

TENTATIVELY IDENTIFIED COMPOUNDS

		Match	ГОР	Amount
Compound	CAS#			hddd
Cyclohexene	110-83-8	NA		Not Detected
Epichlorohydrin	106-89-8	ΑN		Not Detected
Ethyl Acetate	141-78-6	ΑN		Not Detected
Isobutanol	78-83-1	ΑN		Not Detected
Isopropyl ether	108-20-3	Ϋ́		Not Detected
Methacrylonitrile	126-98-7	ΑN		Not Detected
Methyl Acetate	79-20-9	ΑN		Not Detected
N,N-Dimethyl Aniline	121-69-7	ΑN		Not Detected
n-Butanol	71-36-3	Ϋ́		Not Detected
Nitrobenzene	98-95-3	ΑN		Not Detected
Pentane	109-66-0	Ϋ́		Not Detected
Propylene	115-07-1	72%		2.1 NJ
Pyridine	110-86-1	Ϋ́		Not Detected

E = Exceeds instrument calibration range.
J = Estimated value.
NJ =The identification is based on presumptive evidence; estimated value.
D: Analyte not within the DoD scope of accreditation.

Surrogates	CAS#	Limits	%Recovery
1,2-Dichloroethane-d4	17060-07-0	70-130	86
4-Bromofluorobenzene	460-00-4	70-130	26
Toluene-d8	2037-26-5	70-130	96





Air Toxics

MODIFIED EPA METHOD TO-15 GC/MS SIM/FULL SCAN BMS Bldg 3

1.45 msdv.i / v050309sim 5/3/21 01:01 PM Instrument/Filename: Date/Time Analyzed: Dilution Factor: 6 Liter Summa Canister (SIM Certified) Date/Time Collected: 4/29/21 04:10 PM 2104655-01B Client ID: Lab ID: Media:

Compound CAS# (ug/m3) 1,1,1-Trichloroethane 71-55-6 0.045 0.14 1,1,2-Tetrachloroethane 79-34-5 0.11 0.18 1,1,2-Trichloroethane 79-34-5 0.01 0.14 1,1,1-Dichloroethane 75-34-3 0.035 0.10 1,1-Dichloroethane 75-34-3 0.035 0.05 1,2-Dichloroethane 106-93-4 0.049 0.05 1,2-Dichloroethane 106-93-4 0.048 0.05 1,2-Dichloroethane 106-46-7 0.048 0.01 1,2-Dichloroethane 106-46-7 0.048 0.01 Carbon Tetrachloride 56-23-5 0.031 0.06 Chloroethane 75-03-3 0.041 0.01 Chloromethane 75-03-3 0.042 0.01 Chloromethane 76-32-5 0.032 0.01 Chloromethane 76-42-2 0.032 0.01 Ethyl Benzene 76-42-2 0.042 0.18 Freon 12 75-71-8 0.054	(ug/m3) 0.14 0.18 0.10 0.053 0.20 0.10 0.16 0.083 0.069 0.13	(ug/m3) (ug/m3) 0.16 Not Detected 0.20 Not Detected 0.12 Not Detected 0.057 Not Detected 0.057 Not Detected 0.12 Not Detected 0.12 Not Detected 0.12 0.051 J 0.17 Not Detected 0.23 0.46 0.18 0.33 0.19 0.10 J 0.14 0.15 1.1 1.1
71-55-6 0.045 79-34-5 0.11 79-00-5 0.061 79-00-5 0.061 75-34-3 0.035 75-35-4 0.049 106-93-4 0.048 106-46-7 0.048 106-46-7 0.031 56-23-5 0.031 71-43-2 0.031 71-43-2 0.031 71-43-2 0.031 74-87-3 0.042 156-59-2 0.038 100-41-4 0.042 76-14-2 0.057 108-38-3 0.093 91-20-3 0.034	0.14 0.18 0.10 0.053 0.20 0.16 0.083 0.16 0.069 0.069	
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79-00-5 75-34-3 0.035 75-34-3 0.035 75-35-4 0.049 107-06-2 0.048 106-46-7 0.048 106-46-7 0.031 56-23-5 0.031 75-00-3 0.061 74-87-3 0.042 166-59-2 0.038 100-41-4 0.042 76-14-2 0.057 108-38-3 0.041 1634-04-4 0.029 91-20-3 95-47-6 0.035	0.14 0.053 0.20 0.10 0.16 0.083 0.16 0.069	
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75-35-4 0.049 106-93-4 0.088 107-06-2 0.048 106-46-7 0.048 71-43-2 0.01 56-23-5 0.01 75-00-3 0.030 67-66-3 0.061 74-87-3 0.042 100-41-4 0.038 76-14-2 0.054 75-71-8 0.057 108-38-3 0.041 1634-04-4 0.029 91-20-3 0.034	0.053 0.20 0.10 0.16 0.083 0.069 0.054	
DB) 106-93-4 0.088 107-06-2 0.048 106-46-7 0.016 71-43-2 0.031 56-23-5 0.031 75-00-3 0.030 67-66-3 0.061 74-87-3 0.042 100-41-4 0.038 76-14-2 0.054 75-71-8 0.057 108-38-3 0.041 1634-04-4 0.029 91-20-3 0.034	0.20 0.10 0.16 0.083 0.16 0.069 0.13	
107-06-2 0.048 106-46-7 0.16 71-43-2 0.031 56-23-5 0.11 75-00-3 0.030 67-66-3 0.061 74-87-3 0.061 166-59-2 0.038 100-41-4 0.042 76-14-2 0.054 75-71-8 0.057 108-38-3 0.041 1634-04-4 0.029 91-20-3 0.034	0.10 0.16 0.083 0.16 0.069 0.054	
106-46-7 0.16 71-43-2 0.031 56-23-5 0.11 75-00-3 0.030 67-66-3 0.061 74-87-3 0.042 156-59-2 0.038 100-41-4 0.042 76-14-2 0.054 75-71-8 0.057 108-38-3 0.041 1634-04-4 0.029 91-20-3 0.034	0.16 0.083 0.16 0.069 0.13	
71-43-20.03156-23-50.1175-00-30.03067-66-30.06174-87-30.042156-59-20.038100-41-40.04276-14-20.05475-71-80.057108-38-30.0411634-04-40.02991-20-30.034	0.083 0.16 0.069 0.13	
56-23-50.1175-00-30.03067-66-30.06174-87-30.042156-59-20.038100-41-40.04276-14-20.05475-71-80.057108-38-30.0411634-04-40.02991-20-30.034	0.16 0.069 0.13 0.054	
75-00-30.03067-66-30.06174-87-30.042156-59-20.038100-41-40.04276-14-20.05475-71-80.057108-38-30.0411634-04-40.02991-20-30.034	0.069 0.13 0.054	
67-66-30.06174-87-30.042156-59-20.038100-41-40.04276-14-20.05475-71-80.057108-38-30.0411634-04-40.02991-20-30.034	0.13	
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76-14-2 0.054 75-71-8 0.057 108-38-3 0.041 1634-04-4 0.029 9 91-20-3 0.093	0.11	
75-71-8 0.057 108-38-3 0.041 1634-04-4 0.029 91-20-3 0.093 95-47-6 0.034	0.18	
108-38-3 0.041 1634-04-4 0.029 9 91-20-3 0.093 95-47-6 0.034	0.13	3.6
outyl ether 1634-04-4 0.029 91-20-3 0.093 95-47-6 0.034	0.11	
ene 91-20-3 0.093 95-47-6 0.034	0.094	0.52 Not Del
95-47-6 0.034	0.19	0.38 0.22 J
	0.11	
Tetrachloroethene 127-18-4 0.080 0.18	0.18	0.20 0.16 J
Toluene 108-88-3 0.052 0.098	0.098	
trans-1,2-Dichloroethene 156-60-5 0.044 0.10		0.57 Not Detected

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MODIFIED EPA METHOD TO-15 GC/MS SIM/FULL SCAN BMS Bldg 3

2104655-01B Client ID: Lab ID:

Date/Time Collected: 4/29/21 04:10 PM

6 Liter Summa Canister (SIM Certified)

Media:

5/3/21 01:01 PM Date/Time Analyzed: Dilution Factor:

1.45 msdv.i / v050309sim Instrument/Filename:

		MDL	ГОБ	Rpt. Limit	Amount
Compound	CAS#	(ng/m3)	(ng/m3)	(ng/m3)	(ng/m3)
Trichloroethene	79-01-6	090'0	0.14	0.16	0.28
Vinyl Chloride	75-01-4	0.020	0.034	0.037	Not Detected

J = Estimated value. D: Analyte not within the DoD scope of accreditation.

Surrogates	CAS#	Limits	%Recovery
1,2-Dichloroethane-d4	17060-07-0	70-130	86
4-Bromofluorobenzene	460-00-4	70-130	96
Toluene-d8	2037-26-5	70-130	86

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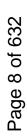
For Laboratory Use Only

180 Blue Ravine Rd. Suite B, Folsom, CA 95630 Air Toxics PID Workorder#

Click links below to view: Canister Sampling Guide

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with all applicable local, State, Federal, and international laws, regulations, and Toxics against any claim, demand, or action, of any kind, related to the collection	eral, and internation or action, of any kin	State, Fede	e local, s y claim,	applicable	oliance with all ns Air Toxics	ipped in compermental in compermental in compermental in compermental in compermental in compermental in compe	ples are sh	les that sam rmless, defe	ment indicat	nature on this docu ndicates agreemer	Sample Transportation Notice: Relinquishing signature on this document indicates that samples are shipped in compliance with all applicable local, State, Federal, and international laws, regulations, and ordinances of any kind. Relinquishing signature also indicates agreement to hold harmless, defend, and indemnify Eurofins Air Toxics against any claim, demand, or action, of any kind, related to the collection	Samp ordinan
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	Helium Shroud Video	Helium S									hone (800) 985-5955; Fax (916) 351-8279	hone





MODIFIED EPA METHOD TO-15 GC/MS SIM/FULL SCAN BMS Syracuse

Client ID: IA-3 Lab ID: 2105127-01A Date/Time Collected: 5/6/21 04:32 PM Media:		Date/Time Analyzed: Dilution Factor: Instrument/Filename:		5/10/21 04:52 PM 6.65 msd20.i / 20051014		
		MDL	COD	Rpt. Limit	Amount	
Compound	CAS#	(ng/m3)	(ng/m3)	(ng/m3)	(ng/m3)	
1,2,4-Trichlorobenzene	120-82-1	3.5	15	25	Not Detected	
1,2,4-Trimethylbenzene	95-63-6	0.23	2.0	3.3	Not Detected	
1,2-Dichlorobenzene	95-50-1	0.39	2.4	4.0	Not Detected	
1,2-Dichloropropane	78-87-5	0.90	1.8	3.1	Not Detected	
1,3,5-Trimethylbenzene	108-67-8	0.46	2.0	3.3	Not Detected	
1,3-Butadiene	106-99-0	0.30	0.88	1.5	Not Detected	
1,3-Dichlorobenzene	541-73-1	0.43	2.4	4.0	Not Detected	
1,4-Dioxane	123-91-1	0.50	4.1	2.4	Not Detected	
2,2,4-Trimethylpentane	540-84-1	1.6	9.3	16	Not Detected	
2-Butanone (Methyl Ethyl Ketone)	78-93-3	1.1	5.9	9.8	Not Detected	
2-Hexanone	591-78-6	1.2	8.2	14	Not Detected	
2-Propanol	67-63-0	1.9	4.9	8.2	89	
3-Chloropropene	107-05-1	1.6	6.2	10	Not Detected	
4-Ethyltoluene	622-96-8	0.39	2.0	3.3	Not Detected	
4-Methyl-2-pentanone	108-10-1	0.33	1.6	2.7	Not Detected	
Acetone	67-64-1	7.0	6.7	16	11 J	
alpha-Chlorotoluene	100-44-7	0.70	2.1	3.4	Not Detected	
Bromodichloromethane	75-27-4	0.63	2.7	4.4	Not Detected	
Bromoform	75-25-2	0.75	4.1	6.9	Not Detected	
Bromomethane	74-83-9	5.9	12	13	Not Detected J	
Carbon Disulfide	75-15-0	0.92	6.2	10	Not Detected	
Chlorobenzene	108-90-7	0.44	1.8	3.1	Not Detected	
cis-1,3-Dichloropropene	10061-01-5	0.49	1.8	3.0	Not Detected	
Cumene	98-82-8	0.26	2.0	3.3	Not Detected	

Page 5 of 26





MODIFIED EPA METHOD TO-15 GC/MS SIM/FULL SCAN BMS Syracuse

6.65 msd20.i / 20051014 5/10/21 04:52 PM Instrument/Filename: Date/Time Analyzed: Dilution Factor: **Lab ID:** 2105127-01A **Date/Time Collected:** 5/6/21 04:32 PM Client ID: Media:

		MDL	ГОР	Rpt. Limit	Amount
Compound	CAS#	(ng/m3)	(ng/m3)	(ng/m3)	(ng/m3)
Cyclohexane	110-82-7	1.8	6.9	11	Not Detected
Dibromochloromethane	124-48-1	0.59	3.4	5.7	Not Detected
Ethanol	64-17-5	2.3	3.8	6.3	1600 E J
Freon 11	75-69-4	1.0	2.2	3.7	1.1 J
Freon 113	76-13-1	0.75	3.0	5.1	Not Detected
Heptane	142-82-5	1.4	8.2	14	Not Detected
Hexachlorobutadiene	87-68-3	5.0	21	35	Not Detected
Hexane	110-54-3	1.1	7.0	12	Not Detected
Methylene Chloride	75-09-2	2.2	4.2	4.6	Not Detected
Propylbenzene	103-65-1	0.32	2.0	3.3	Not Detected
Styrene	100-42-5	0.29	1.7	2.8	Not Detected
Tetrahydrofuran	109-99-9	1.1	5.9	9.8	Not Detected
trans-1,3-Dichloropropene	10061-02-6	0.57	1.8	3.0	Not Detected

D: Analyte not within the DoD scope of accreditation.

TENTATIVELY IDENTIFIED COMPOUNDS

		Match	ТОР	Amount
Compound	CAS#			ppbv
1,2,3-Trimethylbenzene	526-73-8	AN		Not Detected
2-Nitropropane	79-46-9	Ϋ́		Not Detected
4-Chlorotoluene	106-43-4	ΝΑ		Not Detected
Acetonitrile	75-05-8	ΝΑ		Not Detected
Benzaldehyde	100-52-7	Ϋ́		Not Detected
bis(2-Chloroethyl) Ether	111-44-4	Ϋ́		Not Detected





MODIFIED EPA METHOD TO-15 GC/MS SIM/FULL SCAN **BMS Syracuse**

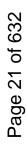
6.65 msd20.i / 20051014 5/10/21 04:52 PM Instrument/Filename: Date/Time Analyzed: Dilution Factor: 6 Liter Summa Canister (SIM Certified) Date/Time Collected: 5/6/21 04:32 PM 2105127-01A Client ID: Lab ID: Media:

TENTATIVELY IDENTIFIED COMPOUNDS

		Match	COD	Amount
Compound	CAS#			h
Cyclohexene	110-83-8	NA		Not Detected
Epichlorohydrin	106-89-8	ΝΑ		Not Detected
Ethyl Acetate	141-78-6	ΝΑ		Not Detected
Isobutanol	78-83-1	ΝΑ		Not Detected
Isopropyl ether	108-20-3	ΝΑ		Not Detected
Methacrylonitrile	126-98-7	ΝΑ		Not Detected
Methyl Acetate	79-20-9	ΝΑ		Not Detected
N,N-Dimethyl Aniline	121-69-7	ΝΑ		Not Detected
n-Butanol	71-36-3	ΝΑ		Not Detected
Nitrobenzene	98-95-3	ΝΑ		Not Detected
Pentane	109-66-0	ΥN		Not Detected
Propylene	115-07-1	ΥN		Not Detected
Pyridine	110-86-1	Ϋ́		Not Detected
1 1 2 4 3 2 4 2 4 2 4 4 4 4 4 4 4 4 4 4 4 4				

J = Estimated value.
E = Exceeds instrument calibration range.
D: Analyte not within the DoD scope of accreditation.

Surrogates	CAS#	Limits	%Recovery
1,2-Dichloroethane-d4	17060-07-0	70-130	104
4-Bromofluorobenzene	460-00-4	70-130	102
Toluene-d8	2037-26-5	70-130	100





MODIFIED EPA METHOD TO-15 GC/MS SIM/FULL SCAN BMS Syracuse

6.65 msd20.i / 20051014sim 5/10/21 04:52 PM Instrument/Filename: Date/Time Analyzed: Dilution Factor: 6 Liter Summa Canister (SIM Certified)
 Lab ID:
 2105127-01B

 Date/Time Collected:
 5/6/21 04:32 PM
 Client ID: Media:

		MDL	ТОР	Rpt. Limit	Amount
Compound	CAS#	(ng/m3)	(ng/m3)	(ng/m3)	(ng/m3)
1,1,1-Trichloroethane	71-55-6	0.041	69:0	0.72	Not Detected
1,1,2,2-Tetrachloroethane	79-34-5	09:0	0.87	0.91	Not Detected
1,1,2-Trichloroethane	79-00-5	0.14	69.0	0.72	Not Detected
1,1-Dichloroethane	75-34-3	0.075	0.51	0.54	Not Detected
1,1-Dichloroethene	75-35-4	0.016	0.053	0.26	Not Detected
1,2-Dibromoethane (EDB)	106-93-4	0.29	0.97	1.0	Not Detected
1,2-Dichloroethane	107-06-2	960.0	0.51	0.54	Not Detected
1,4-Dichlorobenzene	106-46-7	0.52	0.76	0.80	Not Detected
Benzene	71-43-2	0.087	0.40	1.1	0.21 J
Carbon Tetrachloride	56-23-5	0.21	08.0	0.84	0.28 J
Chloroethane	75-00-3	0.063	0.33	0.88	Not Detected
Chloroform	67-66-3	0.14	0.62	0.65	0.15 J
Chloromethane	74-87-3	0.11	0.26	6.9	U.87 J
cis-1,2-Dichloroethene	156-59-2	0.082	0:20	0.53	Not Detected
Ethyl Benzene	100-41-4	0.11	0.55	0.58	Not Detected
Freon 114	76-14-2	0.12	0.88	0.93	Not Detected
Freon 12	75-71-8	0.13	0.62	16	2.0 J
m,p-Xylene	108-38-3	0.12	0.55	1.2	0.28 J
Methyl tert-butyl ether	1634-04-4	0.12	0.46	2.4	Not Detected
Naphthalene	91-20-3	0.23	0.26	1.7	Not Detected
o-Xylene	95-47-6	0.11	0.55	0.58	Not Detected
Tetrachloroethene	127-18-4	0.20	98.0	0.90	Not Detected
Toluene	108-88-3	0.079	0.48	1.2	0.72 J
trans-1,2-Dichloroethene	156-60-5	0.059	0.50	2.6	Not Detected

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MODIFIED EPA METHOD TO-15 GC/MS SIM/FULL SCAN

BMS Syracuse

6.65 msd20.i / 20051014sim 5/10/21 04:52 PM Instrument/Filename: Date/Time Analyzed: Dilution Factor: 6 Liter Summa Canister (SIM Certified)
 Lab ID:
 2105127-01B

 Date/Time Collected:
 5/6/21 04:32 PM
 Client ID: Media:

		MDL	TOD	Rpt. Limit	Amount
Compound	CAS#	(ng/m3)	(ng/m3)	(ng/m3)	(ng/m3)
Trichloroethene	79-01-6	0.12	0.68	0.71	0.26 J
Vinyl Chloride	75-01-4	0.014	0.034	0.17	Not Detected

J = Estimated value. D: Analyte not within the DoD scope of accreditation.

Surrogates	CAS#	Limits	%Recovery
1,2-Dichloroethane-d4	17060-07-0	70-130	66
4-Bromofluorobenzene	460-00-4	70-130	95
Toluene-d8	2037-26-5	70-130	100

STUROUNS

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Phone (800) 985-5955; Fax (916) 351-8279 Shipper Name: Relinquished by: (Signature/Affiliation) Relinquished by: (Signature/Affiliation) Relipiquished by Site Name: Project Manager: Project Name: 180 Blue Ravine Rd. Suite B, Folsom, CA 95630 ordinances of any kind. Relinquishing signature also indicates agreement to hold harmless, defend, and indemnify Eurofins Air Toxics against any claim, demand, or action, of any kind, related to the collection Sample Transportation Notice: Relinquishing signature on this document indicates that samples are shipped in compliance with all applicable local, State, Federal, and international laws, regulations, and Frdex (Signatyre/Affiliatipn Sample Identification BMS SWILLSE 4riadis)an/e Huddis Air Toxics P.O.# Custody Seals Intact? 660453 Can # PIO Date Date handling, of shipping of samples. D.O.T Hotline (800) 467-4922 Controller # 23235 Workorder #: For Laboratory Use Only Special Instructions/Notes: Yes Please Reput Results Date Start Sampling Time Information Time Lab Use Only 0 8 0815 Time None Received by: (Signature/Affiliation) Received by: (Signature/Affiliation) Received by: (Signature/Affiliation) Date Stop Sampling Information 212 Time アカフ Select TAT from drop down box Canister Vacuum/Pressure Initial (in Hg) Turnaround Time (Rush surcharges may apply) Click links below to view: Helium Shroud Video Final (in Hg) Lab Use Only Receipt Date Date 05/14/20 Final (psig) Gas: N₂ / He BMS LO 1015 List 1TTG Requested Analyses me Time Ime 500]





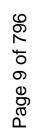
MODIFIED EPA METHOD TO-15 GC/MS SIM/FULL SCAN BMS Syracuse

Lab ID: 2105279A-01A **Date/Time Collected:** 5/13/21 04:15 PM Client ID: Lab ID:

5/17/21 07:24 PM Date/Time Analyzed: Dilution Factor:

Media:		Instrument/Filename:		14.3 msdv.i / v051717		
		4	2	- + c Q	tail out	_
Compound	#S\D	(ug/m3)	(ua/m3)	(ug/m3)	(ng/m3)	
1,2,4-Trichlorobenzene	120-82-1	24	42	53	Not Detected	_
1,2,4-Trimethylbenzene	95-63-6	3.0	5.6	7.0	Not Detected	
1,2-Dichlorobenzene	95-50-1	5.1	6.9	8.6	Not Detected	
1,2-Dichloropropane	78-87-5	3.1	5.3	9.9	Not Detected	
1,3,5-Trimethylbenzene	108-67-8	3.6	5.6	7.0	Not Detected	
1,3-Butadiene	106-99-0	1.3	2.5	3.2	Not Detected	
1,3-Dichlorobenzene	541-73-1	2.6	6.9	8.6	Not Detected	
1,4-Dioxane	123-91-1	2.9	4.1	5.2	Not Detected	
2,2,4-Trimethylpentane	540-84-1	7.4	27	33	Not Detected	
2-Butanone (Methyl Ethyl Ketone)	78-93-3	3.4	17	21	Not Detected	
2-Hexanone	591-78-6	4.2	23	29	Not Detected	
2-Propanol	67-63-0	3.8	14	18	82	
3-Chloropropene	107-05-1	11	18	22	Not Detected	
4-Ethyltoluene	622-96-8	2.7	5.6	7.0	Not Detected	
4-Methyl-2-pentanone	108-10-1	1.3	4.7	5.8	Not Detected	
Acetone	67-64-1	6.1	14	34	11 J	
alpha-Chlorotoluene	100-44-7	4.6	5.9	7.4	Not Detected	
Bromodichloromethane	75-27-4	3.0	7.7	9.6	Not Detected	
Bromoform	75-25-2	4.6	12	15	Not Detected	
Bromomethane	74-83-9	5.4	22	28	Not Detected J	
Carbon Disulfide	75-15-0	3.3	18	22	Not Detected	
Chlorobenzene	108-90-7	2.0	5.3	9.9	Not Detected	
cis-1,3-Dichloropropene	10061-01-5	2.5	5.2	6.5	Not Detected	
Cumene	98-82-8	0.92	5.6	7.0	Not Detected	

Page 5 of 27





MODIFIED EPA METHOD TO-15 GC/MS SIM/FULL SCAN BMS Syracuse

Client ID:

Media:

Lab ID: 2105279A-01A **Date/Time Collected:** 5/13/21 04:15 PM

5/17/21 07:24 PM Date/Time Analyzed:

14.3 msdv.i / v051717 Instrument/Filename: Dilution Factor:

		MDL	ГОР	Rpt. Limit	Amount
Compound	CAS#	(ng/m3)	(ng/m3)	(ng/m3)	(ng/m3)
Cyclohexane	110-82-7	1.9	3.9	25	Not Detected
Dibromochloromethane	124-48-1	4.8	9.7	12	Not Detected
Ethanol	64-17-5	3.9	11	13	3300 E J
Freon 11	75-69-4	1.8	6.4	8.0	Not Detected
Freon 113	76-13-1	4.5	8.8	11	Not Detected
Heptane	142-82-5	5.4	23	29	Not Detected
Hexachlorobutadiene	87-68-3	24	61	92	Not Detected
Hexane	110-54-3	4.2	20	25	Not Detected
Methylene Chloride	75-09-2	2.4	4.0	9.6	Not Detected
Propylbenzene	103-65-1	4.1	5.6	7.0	Not Detected
Styrene	100-42-5	1.7	4.9	6.1	Not Detected
Tetrahydrofuran	109-99-9	7.2	17	21	Not Detected
trans-1,3-Dichloropropene	10061-02-6	2.7	5.2	6.5	Not Detected

D: Analyte not within the DoD scope of accreditation.

TENTATIVELY IDENTIFIED COMPOUNDS

		Match	ГОР	Amount
Compound	CAS#			ppbv
1,2,3-Trimethylbenzene	526-73-8	NA		Not Detected
2-Nitropropane	79-46-9	ΥN		Not Detected
4-Chlorotoluene	106-43-4	ΥN		Not Detected
Acetonitrile	75-05-8	ΥN		Not Detected
Benzaldehyde	100-52-7	ΥN		Not Detected
bis(2-Chloroethyl) Ether	111-44-4	NA		Not Detected





MODIFIED EPA METHOD TO-15 GC/MS SIM/FULL SCAN **BMS Syracuse**

Client ID:

Lab ID: 2105279A-01A **Date/Time Collected:** 5/13/21 04:15 PM

6 Liter Summa Canister (SIM Certified)

Media:

5/17/21 07:24 PM 14.3 msdv.i / v051717 Instrument/Filename: Date/Time Analyzed: Dilution Factor:

TENTATIVELY IDENTIFIED COMPOUNDS

		Match	ГОБ	Amount
Compound	CAS#			hpbv
Cyclohexene	110-83-8	NA		Not Detected
Epichlorohydrin	106-89-8	ΝΑ		Not Detected
Ethyl Acetate	141-78-6	NA		Not Detected
Isobutanol	78-83-1	ΝΑ		Not Detected
Isopropyl ether	108-20-3	NA		Not Detected
Methacrylonitrile	126-98-7	ΝΑ		Not Detected
Methyl Acetate	79-20-9	ΝΑ		Not Detected
N,N-Dimethyl Aniline	121-69-7	NA		Not Detected
n-Butanol	71-36-3	ΝΑ		Not Detected
Nitrobenzene	98-95-3	ΝΑ		Not Detected
Pentane	109-66-0	Ϋ́		Not Detected
Propylene	115-07-1	Ϋ́		Not Detected
Pyridine	110-86-1	Ϋ́		Not Detected

E = Exceeds instrument calibration range.J = Estimated value.D: Analyte not within the DoD scope of accreditation.

	: 0		
Surrogates	CAS#	CIIIIIIS	/orecovery
1,2-Dichloroethane-d4	17060-07-0	70-130	06
4-Bromofluorobenzene	460-00-4	70-130	96
Toluene-d8	2037-26-5	70-130	93





MODIFIED EPA METHOD TO-15 GC/MS SIM/FULL SCAN

BMS Syracuse

IA-3 Client ID: Lab ID

Lab ID: 2105279A-01B Date/Time Collected: 5/13/21 04:15 PM Media: 6 Liter Summa Ca	2105279A-01B 5/13/21 04:15 PM 6 Liter Summa Canister (SIM	ilM Certified)	Date/Time Analyzed: Dilution Factor: Instrument/Filename:	4: 5/17/21 07:24 PM 14.3 1e: msdv.i / v051717sim	24 PM 1717sim	
			MDL	ГОР	Rpt. Limit	Amount
Compound		CAS#	(ng/m3)	(ng/m3)	(ng/m3)	(ng/m3)
1,1,1-Trichloroethane	71	71-55-6	0.45	1.4	1.6	Not Detected
1,1,2,2-Tetrachloroethane		79-34-5	1.1	1.8	2.0	Not Detected
1,1,2-Trichloroethane	62	79-00-5	09:0	1.4	1.6	Not Detected
1,1-Dichloroethane	75	75-34-3	0.35	1.0	1.2	Not Detected
1,1-Dichloroethene	75	75-35-4	0.49	0.52	0.57	Not Detected
1,2-Dibromoethane (EDB)		106-93-4	0.87	2.0	2.2	Not Detected
1,2-Dichloroethane	10	107-06-2	0.47	1.0	1.2	Not Detected
1,4-Dichlorobenzene	10	106-46-7	1.6	1.6	1.7	Not Detected
Benzene	71	71-43-2	0.30	0.82	2.3	Not Detected
Carbon Tetrachloride	99	56-23-5	1.1	1.6	1.8	Not Detected
Chloroethane	75	75-00-3	0.29	0.68	1.9	Not Detected
Chloroform	29	67-66-3	09:0	1.2	1.4	Not Detected
Chloromethane	74	74-87-3	0.41	0.53	15	1.0 J
cis-1,2-Dichloroethene		156-59-2	0.37	1.0	1.1	Not Detected
Ethyl Benzene	10	100-41-4	0.41	1.1	1.2	Not Detected
Freon 114	92	76-14-2	0.54	1.8	2.0	Not Detected
Freon 12	75	75-71-8	0.56	1.3	35	2.3 J
m,p-Xylene	10	108-38-3	0.40	1.1	2.5	Not Detected
Methyl tert-butyl ether		1634-04-4	0.29	0.93	5.2	Not Detected
Naphthalene	91	91-20-3	0.92	1.9	3.7	Not Detected
- ;				•	(

Page 8 of 27

Not Detected Not Detected Not Detected

3.7 1.2 1.9 2.7 5.7

1.9 1.7 0.97

0.34 0.92

0.79

127-18-4 108-88-3 156-60-5

trans-1,2-Dichloroethene

Tetrachloroethene

Toluene

Naphthalene o-Xylene

95-47-6

0.51

Not Detected 0.70 J





MODIFIED EPA METHOD TO-15 GC/MS SIM/FULL SCAN

BMS Syracuse

Client ID: Lab ID:

Media:

msdv.i / v051717sim 5/17/21 07:24 PM Instrument/Filename: Date/Time Analyzed: Dilution Factor: 6 Liter Summa Canister (SIM Certified) Date/Time Collected: 5/13/21 04:15 PM 2105279A-01B

Not Detected Not Detected (ng/m3) Amount Rpt. Limit (ng/m3) 0.36 (ng/m3) LOD 0.34 (ng/m3) 0.59 MDL 0.20 CAS# 79-01-6 75-01-4 **Trichloroethene** Vinyl Chloride Compound

J = Estimated value. D: Analyte not within the DoD scope of accreditation.

Surrogates	CAS#	Limits	%Recovery
1,2-Dichloroethane-d4	17060-07-0	70-130	06
4-Bromofluorobenzene	460-00-4	70-130	26
Toluene-d8	2037-26-5	70-130	26

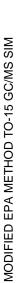
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Analysis Requestive anister Chain of Custody

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180 Blue Ravine Rd. Suite B, Folsom, CA 95630	30 - 50.	vvorkorder #:	Marie Marie	2705773		Click links below to view: Canister Sampling Guide	o view: Guide	
⁹ hone (800) 985-5955; Fax (916) 351-8279					<u> </u>	Helium Shroud Video		
Slient: Arcadis	PID:	Special Ins	Special Instructions/Notes:		Turr	around Time (Ru	Turnaround Time (Rush surcharges may apply)	ay apply)
Project Name: BMS SWaluse			Olar Pan	4 Results	Select TAT	Select TAT from drop down box	oox 3 Day	
Project Manager: Dolling Euch	P.O.#	<u></u>	TROOK RT	•	Caniste	Canister Vacuum/Pressure		Requested Analyses
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Lab Sample Identification	Can #	Flow Controller#	Start Sampling Information	Stop Sampling Information	al (in H	1.1	11 (psig) : N2 / H 113 H Hist	
			Date Time	Date	Time Initia	Rec	Gas B/	
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Sample Transportation Notice: Relinquishing signature on this document indicates that samples are shipped in compliance with all applicable local, State, Federal, and international laws, regulations, an ordinances of any kind. Relinquishing signature also indicates agreement to hold harmless, defend, and indemnify Function Samples agreement or action of any kind. Relinquishing signature also indicates agreement to hold harmless, defend, and indemnify Function Samples agreement or action of any kind.	inature on this docu	ment indicates #	lat samples are shippe is, defend, and indemi	ed in compliance with all a	applicable local, State	e, Federal, and int	ernational laws, rec	gulations, an
		handling, of ship	ping of samples. D.O.	handling, of shipping of samples. D.O.T Hotline (800) 467-4922	Jainst ally Gallit, Gen	iano, or action, or	any kino, selateu to	o the collecti
		and the same of th						





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BMS Syracuse

msd20.i / 20052008sim 5/20/21 12:41 PM Instrument/Filename: Date/Time Analyzed: Dilution Factor: 6 Liter Summa Canister (SIM Certified) **Lab ID:** 2105279B-01A **Date/Time Collected:** 5/13/21 04:15 PM Client ID: Media:

Compound CAS# (ug/m3) (ug/m3)			MDL	TOD	Rpt. Limit		Amount	
79-01-6 0.084 0.48 0.51 0.51 the DoD scope of accreditation. 0.084 0.091 0.091	Compound	CAS#	(ng/m3)	(ng/m3)	(ng/m3)		(ng/m3)	
J = Estimated value. D: Analyte not within the DoD scope of accreditation.	Trichloroethene	79-01-6	0.084	0.48	0.51	0.51	0287	NB
	J = Estimated value. D: Analyte not within the DoD scope	of accreditation.						

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Surrogates	CAS#	Limits	%Recovery
1,2-Dichloroethane-d4	17060-07-0	70-130	102
4-Bromofluorobenzene	460-00-4	70-130	98
Toluene-d8	2037-26-5	70-130	100