#### JSC TOXICOLOGY AND ENVIRONMENTAL CHEMISTRY GROUP

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SUBJECT: Toxicological Assessment of ISS Air and Water Quality: April 10, 2017 – June 2, 2017 (Increment 51), Including OA-7 Ingress Report

SUMMARY: Based on these data, air quality was acceptable on ISS for this period and potable water remains acceptable for crew consumption.

#### **AIR QUALITY**

Three archive air samples were collected in mini grab sample containers (mGSCs) on ISS during Increment 51. Two of these samples were collected as part of routine monitoring in the US Lab and Columbus Module (COL). The third sample was collected as part of nominal ingress operations following initial hatch opening of Orbital/ATK-7 (OA-7). Two pairs of passive-diffusion formaldehyde badges were also deployed in the Lab and SM on 5/8/2017. The mGSC samples were returned on SpX-11 while formaldehyde badges were returned later on SpX-12. A summary of the analytical results from the samples is provided in Table 1.

Sample Location	Sample Date	Freon 218 (mg/m <sup>3</sup> )	Alcohols <sup>a</sup> (mg/m <sup>3</sup> )	T-Value <sup>b</sup> (units)	CO <sub>2</sub> (mg/m <sup>3</sup> )	Formaldehyde (µg/m <sup>3</sup> )
OA-7 Ingress	4/22/2017	21	6.5	2.8 (1.6)	4100	-
Lab	5/8/2017	74	7.5	0.2	6600	25°
COL	5/8/2017	75	7.4	0.2	7000	-
SM	5/8/2017	-	-	-	-	17 <sup>e</sup>
Guideline			<5	$< l^c$	$< 7100^{d}$	<120

Table 1. Analytical summary of ISS air analyses

<sup>a</sup>Includes acetone

<sup>b</sup>Sum of the ratios of the measured concentration and the corresponding 180-day SMAC for each compound, excluding CO<sub>2</sub>; parentheses indicate value based on 7-day SMACs and applicable to first ingress

°T-value <1 used to evaluate routine monthly sampling; <3 used to evaluate first ingress

<sup>d</sup>CO<sub>2</sub> to be controlled as low as reasonably achievable (ALARA) – currently 3 mmHg (7100 mg/m<sup>3</sup>) or lower

<sup>e</sup>Average from pair of formaldehyde badges

Data tables containing measured concentrations and corresponding T-values based on appropriate Spacecraft Maximum Allowable Concentrations (SMACs) for compounds present at levels above the laboratory reporting limit are enclosed. Complete data tables including compounds assessed but not detected are available upon request. The mean relative recoveries of the 3 surrogate standards from the SpX-11 return mGSC samples were as follows: <sup>13</sup>C-acetone, 112±10%; fluorobenzene-d<sub>5</sub>, 108±9%; and chlorobenzene-d<sub>5</sub>, 113±23%. For the passive-diffusion formaldehyde badges, positive control recoveries (1 in-flight and 2 lab controls) were 95, 71, and 106%, respectively.

Automated sampling sessions are scheduled on the Air Quality Monitors (AQMs) every 73 hours, which results in 2-3 sampling sessions per unit per week. Monthly average concentrations as well as the Increment average concentrations for compounds measured on the AQMs are presented in Table 2.

	April	May	Average
2-Propanol	0.19	0.13	0.16
Acetone	0.37	0.35	0.36
Acrolein	ND	ND	ND
Benzene	ND	ND	ND
1,2-Dichloroethane	ND	ND	ND
Decamethylcyclopentasiloxane#	0.16	0.20	0.18
Hexanal	ND	ND	ND
Hexane	ND	ND	ND
m,p-Xylenes#	ND	ND	ND
Methanol	0.29	0.27	0.28
o-Xylene#	0.03	0.03	0.03
Octamethylcylcotetrasiloxane#	TRACE	TRACE	TRACE
Toluene#	0.03	0.03	0.03
2-Butanone	ND	ND	ND
Acetaldehyde	0.12	0.12	0.12
Dichloromethane	ND	ND	ND
Ethanol	3.58	3.71	3.65
Ethyl Acetate	0.05	0.05	0.05
Hexamethycyclotrisiloxane#	0.07	0.07	0.07
n-Butanol	0.08	0.08	0.08
Trimethylsilanol	0.17	0.16	0.17

Table 2. Average monthly concentrations (mg/m<sup>3</sup>) of AQM target compounds

# Obtained from prime unit

ND: Not detected

TRACE= >MDL (Minimum Detection Limit), <MQL (Minimum Quantification Limit)

#### **Toxicological Evaluation of ISS Air Quality**

Routine air quality monitoring is performed in-flight using the AQMs. Archive air samples (mGSCs and formaldehyde badges) are collected during each Increment and returned for analysis in the Toxicology and Environmental Chemistry (TEC) Air Quality Laboratory. Data from the ground analyses complement the in-flight data and provide a more complete understanding of air quality on the ISS. The routine archive samples for this Increment that returned on SpX-11 confirmed air quality was acceptable during this timeframe. All measured values for routine samples (mGSC and AQM) met T-value guideline criteria (T < 1), indicating no concern for crew health. The average, rounded T-value calculated from the Increment 51 mGSC samples was 0.2 (Figure 1). The average, rounded T-value calculated from the AQM data (Figure 2) was slightly lower (0.1 units), but still showed close agreement with the mGSC value. Due to a modification of the process for calculating T-values from AQM data, the T-value for Increment 51 is approximately half of what was documented in recent reports (Refer to TOX-AR-2017-04 for details). Overall, the reported concentrations for the compounds detected are consistent with levels detected since installation of the Node 1 carbon filters in May 2015.

The nominal mGSC samples contained a  $CO_2$  concentration below the Increment limit documented in Chit 14468, which requests that the 24 hour average concentration not exceed 3.0 mmHg (7100 mg/m<sup>3</sup>). While mGSC  $CO_2$  sampling provides a snap-shot of the  $CO_2$  concentration, the major constituent analyzer (MCA) routinely monitors  $CO_2$  levels in the US segment. For this reason, data from the MCA is better suited for evaluation of short and long-term trends in  $CO_2$ . The MCA data concentrations fluctuate as a result of multiple factors including the number of crew on ISS, current scrubbing capability, and processes and activities that generate  $CO_2$ . The average 24 hr  $CO_2$  concentration was approximately 3.0 mmHg or less during all stages (3, 5, and 6-crew operations) of the Increment. Brief excursions above the 3.0 mmHg Increment limit were due to MetOx regeneration activities on April 13 and May 22.  $CO_2$  concentrations were, overall, well controlled.







Figure 2. AQM T-values

Alcohol values in May routine samples continued to exceed the guideline of <5 mg/m3, which is intended to protect the water recovery system from risk of overloading. These levels are primarily due to ethanol in the ISS atmosphere. AQM results for ethanol were lower than the levels measured in the mGSCs, with an Increment average of 3.65 mg/m<sup>3</sup>. This difference may be due to temporal and spatial differences between the AQMs and mGSC sample points. Importantly, ethanol levels during the entire Increment did not present a risk for crew health. Formaldehyde levels in the US Lab (shown in Table 1 and

Figure 3) are generally consistent with historic levels and remain below the SMAC of 120  $\mu$ g/m<sup>3</sup>. Nonmethane volatile organic compounds (NMVOCs) were detected at total concentrations of 8.7 and 8.9 mg/m<sup>3</sup> in nominal archive US Lab and Columbus samples, respectively.



Figure 3. Formaldehyde trending in ISS air.

#### **OA-7** Ingress

A first ingress sample was collected on 4/22/2017, approximately three minutes after hatch opening. The concentration of Freon 218 (octafluoropropane), a marker for ISS air dilution of first entry samples, indicated ~ 27% mixing with the ISS atmosphere prior to sample collection. The total T-value (minus CO<sub>2</sub>) was 1.6, which was well below the limit of 3.0 units. After accounting for dilution, exposure to OA-7 vehicle air would have posed no risk to crew health. Comparatively, the level was similar to OA-6, but the Freon 218 concentration suggests that less dilution with ISS air occurred in the OA-6 sample (~11%). This would imply that the OA-6 first ingress environment contained a lower total concentration of volatile contaminants that significantly contributed to the overall T-value than OA-7. No vehicle off-gas test was performed for OA-7, but previous Orbital/ATK vehicles were tested, including OA-6 and OA-4. First ingress of OA-7 and OA-6 yielded similar levels of fluorotrimethylsilane, a compound that significantly impacted the T-value in OA-4. Acetaldehyde (1.0 mg/m<sup>3</sup>), trimethylsilanol (2.1 mg/m<sup>3</sup>), dodecafluoropentane (55 mg/m<sup>3</sup>), fluorotrimethylsilane (0.49 mg/m<sup>3</sup>), and carbon monoxide (8.9 mg/m<sup>3</sup>) were the primary contributors to the OA-7 T-value. The concentration of total NMVOCs was 78 mg/m<sup>3</sup>. As might be expected, the total volatile organic compound (VOC) concentration was higher in the OA-7 first entry sample than in the nominal ISS samples. VOC totals generally have limited toxicological applications in situations where individual VOCs can be quantified, although the total load may have usefulness to ECLSS or other stakeholders. Previous memos have cited a guideline value of 25 mg/m<sup>3</sup> as a general screening indicator for situations where totals are sufficient to potentially cause odors or perception of poor air quality. However, NASA JSC Toxicology is discontinuing this comparison to total VOC concentration, as there is not sufficient scientific evidence to suggest that this is an appropriate guideline or consideration in spaceflight applications. NASA JSC Toxicology will continue to focus on the evaluation of each measured chemical by comparing with applicable SMACs, and by utilizing t-values to assess the overall toxicity of a chemical mixture.

#### WATER QUALITY

Three archive samples were collected from the US segment during Increment 51. These consisted of one potable water sample from the Hot leg of the US Potable Water Dispenser (PWD), as well as samples of US condensate and wastewater. All three samples were returned on SpX-11. Complete data tables with results for all measured parameters can be found in report 2017-TEC-WQ-003.1. A summary of select analytical results is provided in Tables 3 and 4. Expanded summary tables containing organic carbon recoveries and results for all analytes present at concentrations above reporting limits are included as attachments to this report.

Sample Location	Sample Date	TOC (mg/L)	DMSD (mg/L)	Conductivity (µS/cm)	Total Iodine (mg/L)
PWD (Hot)	4/25/2017	1.39	4.3	2	< 0.05
US Condensate	5/8/2017	48.1	37.0	300	NA
WPA Wastewater	5/19/2017	41.4	12.0	112	NA

Table 3. Analytical Summary of ISS Water Analyses

**Toxicological Evaluation of ISS Water Quality:** Routine water quality monitoring is performed in-flight using the total organic carbon analyzer (TOCA). Results from these analyses provide a general indication of overall water quality. Archive water samples are collected during each Increment and returned for comprehensive analysis in ground laboratories. Data from the ground analyses complement the in-flight data and provide a more complete understanding of water quality on the ISS.

#### Potable Water

Concentrations of all chemicals met the requirements listed in SSP 41000, *System Specification for the International Space Station*. Total organic carbon (TOC) concentrations from in-flight (PWD TOC and WPA PFU2) and ground analyses (Archive TOC) performed on samples from the U.S. potable water system between June 2016 and June 2017 are shown in Figure 4. While the TOC concentration was elevated during Increment 51 compared to normal levels (primarily due to DMSD), the TOC concentrations measured in the U.S. potable water samples and product water sample remained below the U.S. Segment Specification (3000  $\mu$ g/L).



Figure 4. Total Organic Carbon (TOC) trending in US Potable Water

The TOC concentration in the U.S. archive sample (Archive TOC) was 1390 µg/L for the PWD hot sample. Compared to the Increment 50 archive sample (PWD hot: 1330 µg/L), TOC concentrations were similar and well below the Spacecraft Water Exposure Guideline (SWEG) of 5.0 mg/L (5000 µg/L).

As mentioned, the source of the TOC in the potable sample was primarily DMSD (4.3 mg/L). Methyl sulfone, another minor contributor to the TOC, was higher (118 µg/L) than the historical average, but consistent with levels from recent samples. Silicon was also detected (1.29 mg/L) at levels typically seen when DMSD is present in the water. Traces of nickel (3  $\mu$ g/L), aluminum (2  $\mu$ g/L), and zinc (3  $\mu$ g/L) were also detected. Importantly, all chemical parameters measured in U.S. potable water samples collected during Increment 50 met the requirements listed in SSP 41000 and the Medical Operations **Requirement Document (MORD).** 

Iodine is a biocide used on the US segment. It is added to the water produced by the Water Processor Assembly (WPA), but removed prior to crew consumption to avoid potential thyroid dysfunction. The total iodine level in the sample collected from the PWD was below the reporting limit (0.05 mg/L), indicating effective removal of iodine in water intended for consumption. For additional information regarding microbial analyses, please see the Increment 51 post-flight report issued by the JSC Environmental Microbiology Laboratory.

#### Condensate

One condensate sample was collected on 5/8/2017. The TOC level in this sample was 48.1 mg/L, which is below the historical average (166 mg/L). This sample contained some of the lowest levels of contaminants ever detected in a condensate sample. For example, ethanol, 2-propanol, and acetate concentrations were below the method reporting limit. Historical averages for these compounds are 50.1 mg/L, 1.08 mg/L, and 46.0 mg/L, respectively. Metals detected in the sample above 0.1 mg/L included zinc (1.18 mg/L) and nickel (0.501 mg/L). Traces of aluminum (10  $\mu$ g/L), silver (18  $\mu$ g/L), and manganese (25  $\mu$ g/L) were also present. These compounds were effectively removed by the WRS as evidenced by the low or undetectable levels in the potable samples.

#### Wastewater

One wastewater sample was collected on 5/19/2017. The TOC level in this sample (41.4 mg/L) was below the historical average of 46.4 mg/L, but higher than recent samples. The DMSD concentration was 12 mg/L. which was slightly lower than Increment 50 (22 mg/L). Other organic compounds detected at levels above 1 mg/L were ethanol (33.6 mg/L), methanol (9.74 mg/L), acetone (7.82 mg/L), propylene glycol (5.82 mg/L), and ethylene glycol (2.02 mg/L). Metals detected above 0.1 mg/L in the samples were zinc (1.15 mg/L) and nickel (0.117 mg/L). Traces of other metals, including aluminum, manganese, and silver were also present. Chromium concentrations have returned to normal levels, which correlates well with the recent decrease in distillate conductivity following installation of the new Distillation Apparatus (DA) in April 2017. As with the condensate samples, all compounds of toxicological interest were effectively cleaned from the samples by the WRS.

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Concurrence by Valerie Ryder, Ph.D., DABT NASA Toxicologist

Date

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Date

Enclosures Table 1: Analytical concentrations of compounds quantified in mGSCs returned on SpX11 Table 2A: T-values corresponding to concentrations in Table 1, based on 180-day SMACs Table 2B: T-values corresponding to OA-7 Ingress concentrations in Table 1, based on 7day and 180-day SMACs

Table 3: Analytical concentrations of compounds quantified in US potable water sample returned on SpX-11

Table 4: Analytical concentrations of compounds quantified in US wastewater and condensate samples returned on SpX-11

## TABLE 1 ANALYTICAL RESULTS OF SPACEX-11 RETURN AIR SAMPLES

1,1,1,2-Tetrafluoroethane (Norflurane)         Perfluoro(2-methylpentane)         Propene         Propane         Carbonyl sulfide (Carbon oxide sulfide)         Chloromethane         Isobutane         Methanol         Acetaldehyde         2-Methyl-1-propene         Butane         Ethanol *         Acetone         Propanal (Propionaldehyde)         2-Propanol (Isopropanol)         Isoprene (2-Methyl-1,3-butadiene)         2-Methyl-2-propanol         Methylene chloride (Dichloromethane)         Carbon disulfide         1-Propanol         Trimethylsilanol         Butanal (Butyraldehyde)	AQ170137           SN2049           OA-7 Ingress           04/22/17 @           16:51 GMT           16:51 GMT           0.10           0.037           0.026           0.10           TRACE           0.23           0.71           1.0           0.74           0.049           2.5           0.70	(mg/M3) AQ170138 SN2061 LAB 05/08/17 @ 14:30 GMT 0.084 <0.025 <0.025 <0.025 <0.025 <0.025 <0.025 <0.025 <0.025 <0.025 <0.025 <0.025 <0.025 <0.025 <0.025 <0.025 <0.025 <0.025 <0.025 <0.025 <0.025 <0.025 <0.025 <0.025 <0.025 <0.025 <0.025 <0.025 <0.025 <0.025 <0.025 <0.025 <0.025 <0.025 <0.025 <0.025 <0.025 <0.025 <0.025 <0.025 <0.025 <0.025 <0.025 <0.025 <0.025 <0.025 <0.025 <0.025 <0.025 <0.025 <0.025 <0.025 <0.025 <0.025 <0.025 <0.025 <0.025 <0.025 <0.025 <0.025 <0.025 <0.025 <0.025 <0.025 <0.025 <0.025 <0.025 <0.025 <0.025 <0.025 <0.025 <0.025 <0.025 <0.025 <0.025 <0.025 <0.025 <0.025 <0.025 <0.025 <0.025 <0.025 <0.025 <0.025 <0.025 <0.025 <0.025 <0.025 <0.025 <0.025 <0.025 <0.025 <0.025 <0.025 <0.025 <0.025 <0.025 <0.025 <0.025 <0.025 <0.025 <0.025 <0.025 <0.028 TRACE	AQ170139 SN2059 Columbus 05/08/17 @ 14:30 GMT 0.081 <0.025 <0.025 <0.025 <0.025 <0.025 <0.025 <0.025 <0.025 <0.025 0.39
Propene Propane Carbonyl sulfide (Carbon oxide sulfide) Chloromethane Isobutane Methanol Acetaldehyde 2-Methyl-1-propene Butane Ethanol * Acetone Propanal (Propionaldehyde) 2-Propanol (Isopropanol) Isoprene (2-Methyl-1,3-butadiene) 2-Methyl-2-propanol Methylene chloride (Dichloromethane) Carbon disulfide 1-Propanol Trimethylsilanol Butanal (Butyraldehyde)	<0.10         0.037         0.026         0.10         TRACE         0.23         0.71         1.0         0.74         0.049         2.5	<0.10 <0.025 <0.025 <0.025 <0.025 <0.025 0.35 0.28	<0.10 <0.025 <0.025 <0.025 <0.025 <0.025 <0.025
Perfluoro(2-methylpentane) Propene Propane Carbonyl sulfide (Carbon oxide sulfide) Chloromethane Isobutane Methanol Acetaldehyde 2-Methyl-1-propene Butane Ethanol * Acetone Propanal (Propionaldehyde) 2-Propanol (Isopropanol) Isoprene (2-Methyl-1,3-butadiene) 2-Methyl-2-propanol Methylene chloride (Dichloromethane) Carbon disulfide 1-Propanol Trimethylsilanol Butanal (Butyraldehyde)	<0.10	<0.10 <0.025 <0.025 <0.025 <0.025 <0.025 0.35 0.28	<0.10 <0.025 <0.025 <0.025 <0.025 <0.025 <0.025
Propene Propane Carbonyl sulfide (Carbon oxide sulfide) Chloromethane Isobutane Methanol Acetaldehyde 2-Methyl-1-propene Butane Ethanol * Acetone Propanal (Propionaldehyde) 2-Propanol (Isopropanol) Isoprene (2-Methyl-1,3-butadiene) 2-Methyl-2-propanol Methylene chloride (Dichloromethane) Carbon disulfide 1-Propanol Trimethylsilanol Butanal (Butyraldehyde)	0.037           0.026           0.10           TRACE           0.23           0.71           1.0           0.74           0.049           2.5	<0.025 <0.025 <0.025 <0.025 <0.025 0.35 0.28	<0.025 <0.025 <0.025 <0.025 <0.025 <0.025
Propane Carbonyl sulfide (Carbon oxide sulfide) Chloromethane Isobutane Methanol Acetaldehyde 2-Methyl-1-propene Butane Ethanol * Acetone Propanal (Propionaldehyde) 2-Propanol (Isopropanol) Isoprene (2-Methyl-1,3-butadiene) 2-Methyl-2-propanol Methylene chloride (Dichloromethane) Carbon disulfide 1-Propanol Trimethylsilanol Butanal (Butyraldehyde)	0.026           0.10           TRACE           0.23           0.71           1.0           0.74           0.049           2.5	<0.025 <0.025 <0.025 <0.025 0.35 0.28	<0.025 <0.025 <0.025 <0.025
Carbonyl sulfide (Carbon oxide sulfide) Chloromethane Isobutane Methanol Acetaldehyde 2-Methyl-1-propene Butane Ethanol * Acetone Propanal (Propionaldehyde) 2-Propanol (Isopropanol) Isoprene (2-Methyl-1,3-butadiene) 2-Methyl-2-propanol Methylene chloride (Dichloromethane) Carbon disulfide 1-Propanol Trimethylsilanol Butanal (Butyraldehyde)	0.10 TRACE 0.23 0.71 1.0 0.74 0.049 <b>2.5</b>	<0.025 <0.025 <0.025 0.35 0.28	<0.025 <0.025 <0.025
Isobutane Methanol Acetaldehyde 2-Methyl-1-propene Butane Ethanol * Acetone Propanal (Propionaldehyde) 2-Propanol (Isopropanol) Isoprene (2-Methyl-1,3-butadiene) 2-Methyl-2-propanol Methylene chloride (Dichloromethane) Carbon disulfide 1-Propanol Trimethylsilanol Butanal (Butyraldehyde)	0.23 0.71 1.0 0.74 0.049 <b>2.5</b>	<0.025 0.35 0.28	< 0.025
Methanol Acetaldehyde 2-Methyl-1-propene Butane Ethanol * Acetone Propanal (Propionaldehyde) 2-Propanol (Isopropanol) Isoprene (2-Methyl-1,3-butadiene) 2-Methyl-2-propanol Methylene chloride (Dichloromethane) Carbon disulfide 1-Propanol Trimethylsilanol Butanal (Butyraldehyde)	0.71 1.0 0.74 0.049 2.5	0.35 0.28	
Acetaldehyde 2-Methyl-1-propene Butane Ethanol * Acetone Propanal (Propionaldehyde) 2-Propanol (Isopropanol) Isoprene (2-Methyl-1,3-butadiene) 2-Methyl-2-propanol Methylene chloride (Dichloromethane) Carbon disulfide 1-Propanol Trimethylsilanol Butanal (Butyraldehyde)	1.0           0.74           0.049           2.5	0.28	0.20
2-Methyl-1-propene Butane Ethanol * Acetone Propanal (Propionaldehyde) 2-Propanol (Isopropanol) Isoprene (2-Methyl-1,3-butadiene) 2-Methyl-2-propanol Methylene chloride (Dichloromethane) Carbon disulfide 1-Propanol Trimethylsilanol Butanal (Butyraldehyde)	0.74 0.049 <b>2.5</b>		
Butane Ethanol * Acetone Propanal (Propionaldehyde) 2-Propanol (Isopropanol) Isoprene (2-Methyl-1,3-butadiene) 2-Methyl-2-propanol Methylene chloride (Dichloromethane) Carbon disulfide 1-Propanol Trimethylsilanol Butanal (Butyraldehyde)	0.049 <b>2.5</b>		0.28 <0.025
Acetone Propanal (Propionaldehyde) 2-Propanol (Isopropanol) Isoprene (2-Methyl-1,3-butadiene) 2-Methyl-2-propanol Methylene chloride (Dichloromethane) Carbon disulfide 1-Propanol Trimethylsilanol Butanal (Butyraldehyde)	2.5	<0.025	< 0.025
Propanal (Propionaldehyde) 2-Propanol (Isopropanol) Isoprene (2-Methyl-1,3-butadiene) 2-Methyl-2-propanol Methylene chloride (Dichloromethane) Carbon disulfide 1-Propanol Trimethylsilanol Butanal (Butyraldehyde)	0.70	6.6	6.4
2-Propanol (Isopropanol) Isoprene (2-Methyl-1,3-butadiene) 2-Methyl-2-propanol Methylene chloride (Dichloromethane) Carbon disulfide 1-Propanol Trimethylsilanol Butanal (Butyraldehyde)		0.31	0.33
Isoprene (2-Methyl-1,3-butadiene) 2-Methyl-2-propanol Methylene chloride (Dichloromethane) Carbon disulfide 1-Propanol Trimethylsilanol Butanal (Butyraldehyde)	0.078	<0.025	< 0.025
2-Methyl-2-propanol Methylene chloride (Dichloromethane) Carbon disulfide 1-Propanol Trimethylsilanol Butanal (Butyraldehyde)	2.3	0.15 0.029	0.15 0.033
Methylene chloride (Dichloromethane) Carbon disulfide 1-Propanol Trimethylsilanol Butanal (Butyraldehyde)	<0.025 0.13	<0.029	<0.033
Carbon disulfide 1-Propanol Trimethylsilanol Butanal (Butyraldehyde)	<0.025	<0.025	<0.025
Trimethylsilanol Butanal (Butyraldehyde)	0.048	<0.025	< 0.025
Butanal (Butyraldehyde)	0.043	0.032	0.035
	2.1	0.14	0.16
2-Butanone (Mathul athul katona)	0.040	<0.025 TRACE	<0.025 TRACE
2-Butanone (Methyl ethyl ketone) Ethyl acetate	0.24	0.028	0.029
1-Butanol	0.17	0.065	0.029
2-Methylhexane	<0.025	<0.025	< 0.025
4-Methyl-2-pentanone (MIBK)	0.026	< 0.025	< 0.025
Toluene	0.11	TRACE	0.027
Hexanal Butul accetate	0.029	<0.025 <0.050	<0.025 <0.050
Butyl acetate Chlorobenzene	<0.050	<0.050	<0.050 TRACE
Heptanal	TRACE	<0.050	<0.050
o-Xylene	TRACE	< 0.050	< 0.050
Octamethylcyclotetrasiloxane	0.99	<0.125	<0.125
1,4-Dichlorobenzene	<0.050	<0.050	0.065
Decamethylcyclopentasiloxane Octafluoropropane (Perfluoropropane) *	0.73 <b>21</b>	0.24 74	0.30 75
SPECIAL INTEREST COMPOUNDS ***		17	15
Hexamethylcyclotrisiloxane #	5.7	<0.20	TRACE
NON-TARGET COMPOUNDS ***		0.051	
Dodecafluoropentane	55	0.081	0.097
Tetradecafluorohexane Fluorotrimethylsilane	0.20 0.49	0.26	0.27 <0.050
2-Methyl-1-propanol	0.060	<0.050	<0.050
Pentamethyldisiloxane-1-ol	0.15	<0.050	<0.050
Octamethyltrisiloxane	0.17	< 0.050	< 0.050
C11-Alkane	0.12	<0.050	< 0.050
Unidentified siloxane C11-Alkane	0.14 0.060	<0.050 <0.050	<0.050 <0.050
C11-Alkane	0.060	<0.050	<0.050
Decamethyltetrasiloxane	0.10	<0.050	<0.050
C12-Alkane	0.071	< 0.050	< 0.050
C12-Alkane	0.23	<0.050	<0.050
C12-Alkane C12-Alkane	0.17 0.17	<0.050 <0.050	<0.050 <0.050
C12-Alkane	0.17	<0.050	<0.050
TOTAL ALCOHOLS PLUS ACETONE	6.5	7.5	7.4
TARGET COMPOUNDS (GC) **			
Methane	4.9	16.2	16.5
Carbon dioxide	4100	6600	7000
Hydrogen	2.2	3.8	3.9
Carbon monoxide	8.9	0.85	0.85
TOTAL CONCENTRATION (NON-METHANE HYDROCARBONS)	99	82	84
TOTAL CONCENTRATION - OFP	78	8.7	8.9

\* GC/FID data results are in bold

\*\* Quantified using a multi-point calibration

\*\*\* Quantified using "B" response factor except where noted; concentrations are estimates only.

# Response factor generated from an internal study

<: Value is less than the laboratory reporting limit.

TRACE: Amount detected is sufficient for compound identification only. One-half of the reporting limit was

used in the Total Concentration summation.

OFP - Octafluoropropane

#### TABLE 2A T-VALUES FOR SPACEX-11 RETURN AIR SAMPLES

CHEMICAL CONTAMINANT		T-VALUE (180-d SMAC)			
	AQ170138 SN2061 LAB 05/08/17 @ 14:30 GMT	AQ170139 SN2059 Columbus 05/08/17 @ 14:30 GMT			
TARGET COMPOUNDS (TO-15)					
1,1,1,2-Tetrafluoroethane (Norflurane)	0.00042	0.00040			
Methanol	0.00393	0.00436			
Acetaldehyde	0.06961	0.07021			
2-Methyl-1-propene	0.00011	ND			
Ethanol	0.00328	0.00318			
Acetone	0.00604	0.00644			
2-Propanol (Isopropanol)	0.00102	0.00098			
Isoprene (2-Methyl-1,3-butadiene)	0.00973	0.01091			
1-Propanol	0.00033	0.00036			
Trimethylsilanol	0.03555	0.04022			
2-Butanone (Methyl ethyl ketone)	0.00042	0.00042			
Ethyl acetate	0.00016	0.00016			
1-Butanol	0.00162	0.00287			
Toluene	0.00083	0.00181			
Chlorobenzene	ND	0.00054			
1,4-Dichlorobenzene	ND	0.00217			
Decamethylcyclopentasiloxane	0.01571	0.02024			
Octafluoropropane (Perfluoropropane)	0.00086	0.00088			
SPECIAL INTEREST COMPOUNDS					
Hexamethylcyclotrisiloxane	ND	0.01111			
NON-TARGET COMPOUNDS					
Dodecafluoropentane	0.00028	0.00033			
Tetradecafluorohexane	0.00000	0.00000			
TARGET COMPOUNDS (GC)					
Methane	0.00464	0.00472			
Carbon dioxide	0.51029	0.53633			
Hydrogen	0.01122	0.01144			
Carbon monoxide	0.04972	0.05026			
Carbon monoxide	0.04772	0.03020			
TOTAL T-VALUE	0.72574	0.78035			
TOTAL T-VALUE - CO2	0.21545	0.24402			

ND : Value is less than the laboratory reporting limit. Note: Number of decimal places in T-Values do not represent significant figures of measurements.

## TABLE 2BT-VALUES FOR OA-7 INGRESS AIR SAMPLE

	T-VALUE (7-d SMAC)	T-VALUE (180-d SMAC)		
CHEMICAL CONTAMINANT	AQ170137 SN2049	AQ170137 SN2049		
	OA-7 Ingress 04/22/17 @ 16:51 GMT	OA-7 Ingress 04/22/17 @ 16:51 GMT		
TARGET COMPOUNDS (TO-15)				
1,1,1,2-Tetrafluoroethane (Norflurane)	0.00631	0.00631		
Propene	0.00021	0.00021		
Propane Carbonyl sulfide (Carbon oxide sulfide)	0.00024 0.04133	0.00469 0.04133		
Chloromethane	0.00030	0.00030		
Isobutane	0.00097	0.00097		
Methanol	0.00787	0.00787		
Acetaldehyde	0.25922	0.25922		
2-Methyl-1-propene	0.00647	0.00647		
Butane	0.00034	0.00693		
Ethanol Acetone	0.00123 0.01347	0.00123 0.01347		
Propanal (Propionaldehyde)	0.00652	0.00652		
2-Propanol (Isopropanol)	0.01540	0.01540		
2-Methyl-2-propanol	0.00090	0.00112		
Carbon disulfide	0.08048	0.08048		
1-Propanol	0.00044	0.00044		
Frimethylsilanol	0.51269	0.51269		
Butanal (Butyraldehyde)	0.00265	0.00265		
2-Butanone (Methyl ethyl ketone)	0.00813	0.00813		
Ethyl acetate 1-Butanol	0.00026	0.00026		
3-Methylhexane	0.00213	0.00431		
4-Methyl-2-pentanone (MIBK)	0.00019	0.00019		
Foluene	0.00727	0.00727		
Hexanal	0.00144	0.00144		
Butyl acetate	0.00058	0.00058		
Heptanal	0.00109	0.00109		
p-Xylene	0.00034	0.00068		
Octamethylcyclotetrasiloxane Decamethylcyclopentasiloxane	0.00355 0.00732	0.08288 0.04879		
Octafluoropropane (Perfluoropropane)				
	0.00024	0.00024		
SPECIAL INTEREST COMPOUNDS	0.00024	0.00024		
SPECIAL INTEREST COMPOUNDS Hexamethylcyclotrisiloxane				
SPECIAL INTEREST COMPOUNDS				
SPECIAL INTEREST COMPOUNDS Hexamethylcyclotrisiloxane NON-TARGET COMPOUNDS Dodecafluoropentane Tetradecafluorohexane	0.06357	0.63572		
SPECIAL INTEREST COMPOUNDS Hexamethylcyclotrisiloxane NON-TARGET COMPOUNDS Dodecafluoropentane Fetradecafluorohexane Fluorotrimethylsilane	0.06357 0.18774 0.00000 0.12832	0.63572 0.18774 0.00000 0.12832		
SPECIAL INTEREST COMPOUNDS Hexamethylcyclotrisiloxane NON-TARGET COMPOUNDS Dodecafluoropentane Fetradecafluorohexane Fluorotrimethylsilane 2-Methyl-1-propanol	0.06357 0.18774 0.00000 0.12832 0.00050	0.63572 0.18774 0.00000 0.12832 0.00050		
SPECIAL INTEREST COMPOUNDS Hexamethylcyclotrisiloxane NON-TARGET COMPOUNDS Dodecafluoropentane Tetradecafluorohexane Fluorotrimethylsilane 2-Methyl-1-propanol Pentamethyldisiloxane-1-ol	0.06357 0.18774 0.00000 0.12832 0.00050 0.00302	0.63572 0.18774 0.00000 0.12832 0.00050 0.00302		
SPECIAL INTEREST COMPOUNDS Hexamethylcyclotrisiloxane NON-TARGET COMPOUNDS Dodecafluoropentane Tetradecafluorohexane Fluorotrimethylsilane 2-Methyl-1-propanol Pentamethyldisiloxane-1-ol Octamethyltrisiloxane	0.06357 0.18774 0.00000 0.12832 0.00050 0.00302 0.00017	0.63572 0.18774 0.00000 0.12832 0.00050 0.00302 0.00425		
SPECIAL INTEREST COMPOUNDS Hexamethylcyclotrisiloxane NON-TARGET COMPOUNDS Dodecafluoropentane Tetradecafluorohexane Fluorotrimethylsilane 2-Methyl-1-propanol Pentamethyldisiloxane-1-ol Octamethyltrisiloxane C11-Alkane	0.06357 0.18774 0.00000 0.12832 0.00050 0.00302 0.00017 0.00265	0.63572 0.18774 0.00000 0.12832 0.00050 0.00302 0.00425 0.00265		
SPECIAL INTEREST COMPOUNDS Hexamethylcyclotrisiloxane NON-TARGET COMPOUNDS Dodecafluoropentane Fetradecafluorohexane Fluorotrimethylsilane 2-Methyl-1-propanol Pentamethyldisiloxane-1-ol Dctamethyltrisiloxane C11-Alkane Unidentified siloxane	0.06357 0.18774 0.00000 0.12832 0.00050 0.00302 0.00017	0.63572 0.18774 0.00000 0.12832 0.00050 0.00302 0.00425		
SPECIAL INTEREST COMPOUNDS Hexamethylcyclotrisiloxane NON-TARGET COMPOUNDS Dodecafluoropentane Tetradecafluorohexane Fluorotrimethylsilane 2-Methyl-1-propanol Pentamethyldisiloxane-1-ol Octamethyltrisiloxane C11-Alkane Unidentified siloxane C11-Alkane	0.06357 0.18774 0.00000 0.12832 0.00050 0.00302 0.00017 0.00265 0.01383	0.63572 0.18774 0.00000 0.12832 0.00050 0.00302 0.00425 0.00265 0.13827		
SPECIAL INTEREST COMPOUNDS Hexamethylcyclotrisiloxane NON-TARGET COMPOUNDS Dodecafluoropentane Tetradecafluorohexane Fluorotrimethylsilane 2-Methyl-1-propanol Pentamethyldisiloxane-1-ol Octamethyltrisiloxane C11-Alkane Unidentified siloxane C11-Alkane C11-Alkane Decamethyltetrasiloxane	0.06357 0.18774 0.00000 0.12832 0.00050 0.00302 0.00017 0.00265 0.01383 0.00136 0.00402 0.00009	0.63572 0.18774 0.00000 0.12832 0.00050 0.00302 0.00425 0.00265 0.13827 0.00136 0.00402 0.00018		
SPECIAL INTEREST COMPOUNDS Hexamethylcyclotrisiloxane NON-TARGET COMPOUNDS Dodecafluoropentane Fetradecafluorohexane Fluorotrimethylsilane 2-Methyl-1-propanol Pentamethyldisiloxane-1-ol Dctamethyltrisiloxane C11-Alkane Unidentified siloxane C11-Alkane C12-Alkane Decamethyltetrasiloxane C12-Alkane	0.06357           0.18774           0.00000           0.12832           0.00050           0.000302           0.00017           0.00265           0.01383           0.00136           0.00009           0.000161	0.63572 0.18774 0.00000 0.12832 0.00050 0.00302 0.00425 0.00265 0.13827 0.00136 0.00402 0.00018 0.00161		
SPECIAL INTEREST COMPOUNDS Hexamethylcyclotrisiloxane NON-TARGET COMPOUNDS Dodecafluoropentane Tetradecafluorohexane Fluorotrimethylsilane 2-Methyl-1-propanol Pentamethyldisiloxane-1-ol Octamethyltrisiloxane C11-Alkane Unidentified siloxane C11-Alkane C11-Alkane C12-Alkane Decamethyltetrasiloxane C12-Alkane C12-Alkane	0.06357           0.18774           0.00000           0.12832           0.00050           0.00302           0.00265           0.01383           0.00136           0.00402           0.0009           0.00161           0.00527	0.63572 0.18774 0.00000 0.12832 0.00050 0.00302 0.00425 0.00265 0.13827 0.00136 0.00402 0.00402 0.00018 0.00161 0.00527		
SPECIAL INTEREST COMPOUNDS Hexamethylcyclotrisiloxane NON-TARGET COMPOUNDS Dodecafluoropentane Tetradecafluorohexane Fluorotrimethylsilane 2-Methyl-1-propanol Pentamethyldisiloxane-1-ol Dctamethyltrisiloxane C11-Alkane Unidentified siloxane C11-Alkane C11-Alkane C12-Alkane Decamethyltetrasiloxane C12-Alkane C12-Alkane C12-Alkane	0.06357           0.18774           0.00000           0.12832           0.00050           0.00302           0.00265           0.01383           0.00136           0.00402           0.00009           0.00161           0.00385	0.63572 0.18774 0.00000 0.12832 0.00050 0.00302 0.00425 0.00265 0.13827 0.00136 0.00402 0.00018 0.00161 0.00527 0.00385		
SPECIAL INTEREST COMPOUNDS Hexamethylcyclotrisiloxane NON-TARGET COMPOUNDS Dodecafluoropentane Fetradecafluorohexane Fluorotrimethylsilane 2-Methyl-1-propanol Pentamethyldisiloxane-1-ol Dctamethyltrisiloxane C11-Alkane C11-Alkane C11-Alkane C11-Alkane C12-Alkane C12-Alkane C12-Alkane C12-Alkane C12-Alkane C12-Alkane C12-Alkane C12-Alkane	0.06357           0.18774           0.00000           0.12832           0.00050           0.00302           0.00265           0.01383           0.00136           0.00402           0.0009           0.00161           0.00527	0.63572 0.18774 0.00000 0.12832 0.00050 0.00302 0.00425 0.00265 0.13827 0.00136 0.00402 0.00018 0.00161 0.00527		
SPECIAL INTEREST COMPOUNDS Hexamethylcyclotrisiloxane NON-TARGET COMPOUNDS Dodecafluoropentane Tetradecafluorohexane Fluorotrimethylsilane 2-Methyl-1-propanol Pentamethyldisiloxane-1-ol Dctamethyltrisiloxane C11-Alkane Unidentified siloxane C11-Alkane C12-Alkane C12-Alkane C12-Alkane C12-Alkane C12-Alkane C12-Alkane C12-Alkane C12-Alkane C12-Alkane	0.06357           0.18774           0.00000           0.12832           0.00050           0.00302           0.00265           0.01383           0.00136           0.00402           0.000527           0.00385           0.00385	0.63572 0.18774 0.00000 0.12832 0.00050 0.00302 0.00425 0.00265 0.13827 0.00136 0.00402 0.00018 0.00161 0.00527 0.00385 0.00385		
SPECIAL INTEREST COMPOUNDS Hexamethylcyclotrisiloxane NON-TARGET COMPOUNDS Dodecafluoropentane Tetradecafluorohexane Fluorotrimethylsilane 2-Methyl-1-propanol Pentamethyldisiloxane-1-ol Octamethyltrisiloxane C11-Alkane Unidentified siloxane C11-Alkane C12-Alkane Decamethyltetrasiloxane C12-Alkane C12-Alkane C12-Alkane C12-Alkane C12-Alkane C12-Alkane C12-Alkane C12-Alkane C12-Alkane C12-Alkane	0.06357           0.18774           0.00000           0.12832           0.00050           0.00302           0.00017           0.00265           0.01383           0.00136           0.000402           0.000161           0.00385           0.00385           0.00143	0.63572 0.18774 0.00000 0.12832 0.00050 0.00302 0.00425 0.00265 0.13827 0.00136 0.00402 0.00018 0.00161 0.00527 0.00385 0.00385 0.00385 0.00143		
SPECIAL INTEREST COMPOUNDS Hexamethylcyclotrisiloxane NON-TARGET COMPOUNDS Dodecafluoropentane Tetradecafluorohexane Fluorotrimethylsilane 2-Methyl-1-propanol Pentamethyldisiloxane-1-ol Octamethyltrisiloxane C11-Alkane Unidentified siloxane C11-Alkane C12-Alkane C12-Alkane C12-Alkane C12-Alkane C12-Alkane C12-Alkane C12-Alkane C12-Alkane C12-Alkane C12-Alkane C12-Alkane C12-Alkane C12-Alkane C12-Alkane C12-Alkane	0.06357           0.18774           0.00000           0.12832           0.00050           0.00050           0.000265           0.01383           0.00136           0.0009           0.00161           0.00385           0.00385           0.00143	0.63572 0.18774 0.00000 0.12832 0.00050 0.00302 0.00425 0.00265 0.13827 0.00136 0.00402 0.00136 0.00161 0.00527 0.00385 0.00385 0.00143 0.00139		
SPECIAL INTEREST COMPOUNDS Hexamethylcyclotrisiloxane NON-TARGET COMPOUNDS Dodecafluoropentane Tetradecafluorohexane Fluorotrimethylsilane 2-Methyl-1-propanol Pentamethyldisiloxane-1-ol Octamethyltrisiloxane C11-Alkane C11-Alkane C11-Alkane C11-Alkane C12-Alkane	0.06357           0.18774           0.00000           0.12832           0.00050           0.00302           0.00017           0.00265           0.00136           0.00402           0.00050           0.00161           0.00385           0.00143           0.00139           0.31446	0.63572           0.18774           0.00000           0.12832           0.00050           0.00302           0.00265           0.13827           0.00136           0.00425           0.00136           0.00018           0.00161           0.00385           0.00143           0.00139           0.31446		
SPECIAL INTEREST COMPOUNDS Hexamethylcyclotrisiloxane NON-TARGET COMPOUNDS Dodecafluoropentane Tetradecafluorohexane Fluorotrimethylsilane 2-Methyl-1-propanol Pentamethyldisiloxane-1-ol Octamethyltrisiloxane C11-Alkane Unidentified siloxane C11-Alkane C12-Alkane Decamethyltetrasiloxane C12-Alkane C12-Alkane C12-Alkane C12-Alkane C12-Alkane C12-Alkane C12-Alkane C12-Alkane C12-Alkane C12-Alkane	0.06357           0.18774           0.00000           0.12832           0.00050           0.00050           0.000265           0.01383           0.00136           0.0009           0.00161           0.00385           0.00385           0.00143	0.63572 0.18774 0.00000 0.12832 0.00050 0.00302 0.00425 0.00265 0.13827 0.00136 0.00402 0.00136 0.00161 0.00527 0.00385 0.00385 0.00143 0.00139		
SPECIAL INTEREST COMPOUNDS Hexamethylcyclotrisiloxane NON-TARGET COMPOUNDS Dodecafluoropentane Fetradecafluorohexane Fluorotrimethylsilane 2-Methyl-1-propanol Pentamethyldisiloxane-1-ol Detamethyltrisiloxane C11-Alkane C11-Alkane C11-Alkane C12-A	0.06357           0.18774           0.00000           0.12832           0.00050           0.00050           0.000265           0.01383           0.00136           0.0009           0.00161           0.00385           0.00385           0.00143	0.63572           0.18774           0.00000           0.12832           0.00050           0.00302           0.00425           0.00136           0.00402           0.00136           0.00161           0.00385           0.00385           0.00143		
SPECIAL INTEREST COMPOUNDS Hexamethylcyclotrisiloxane NON-TARGET COMPOUNDS Dodecafluoropentane Fetradecafluorohexane Fluorotrimethylsilane 2-Methyl-1-propanol Pentamethyldisiloxane-1-ol Octamethyltrisiloxane C11-Alkane Unidentified siloxane C11-Alkane C12-Alkane	0.06357           0.18774           0.00000           0.12832           0.00050           0.00050           0.000265           0.01383           0.00136           0.0009           0.00161           0.00385           0.00385           0.00139           0.31446           0.00650           0.14082	0.63572           0.18774           0.00000           0.12832           0.00050           0.00302           0.00425           0.00136           0.00136           0.00161           0.00385           0.00143           0.00139           0.31446           0.00650           0.52187		

ND : Value is less than the laboratory reporting limit.

Note: Number of decimal places in T-Values do not represent significant figures of measurements.

## Table 3. Increment 51 Water Sample Summary ReportUS Potable Water Sample

Increment					51
Mission					SpX-11
Sample Location			Potable Water		WPA PWD Hot
Sample Description		Test	Maximum Contaminant	Maximum Contaminant	Potable Water
Sample Date		Conducted	Level	Level	4/25/2017
Analysis/Sample ID	Units	by	(MCL)	Source	20170705001
Physical Characteristics					
pH	pH units	U.S.	4.5-8.5	41000	5.87
Conductivity	µS/cm	U.S.			2
Trace Metals (ICP/MS)					
Calcium	mg/L	U.S.	30	41000	0.02
Sodium	mg/L	U.S.			0.03
Aluminum	µg/L	U.S.			2
Nickel	µg/L	U.S.	300	SWEG&41000	3
Zinc	µg/L	U.S.	2,000	SWEG&41000	3
Silicon (ICP/MS)					
Silicon	µg/L	U.S.			1,290
Total Organic Carbon (Sievers)					
Inorganic Carbon	mg/L	U.S.			0.61
Organic Carbon	mg/L	U.S.	3	41000	1.39
Semi-volatiles (GC/MS) - Target List					
Methyl sulfone	μg/L	U.S.	1,500,000	interim SWEG (06-2017)	118
Silanols (LC/RI) (R & D Method -NIST traceable stand	lard not av	ailable)			
Dimethylsilanediol (DMSD)	μg/L	U.S.	35,000	SWEG	4,300
Organic Carbon Recovery	percent	U.S.			82.72
Unaccounted Organic Carbon	mg/L	U.S.			0.24

NA=Not analyzed MI=Matrix Interference N/A=Not applicable

# Table 4. Increment 51 Water Sample Summary Report WPAWastewater and Condensate Samples

Increment						51
Mission					SpX-11	
Sample Location Sample Description		Test	Potable Water Maximum Contaminant	Maximum Contaminant	WPA Wastewater ORU WPA Wastewater	WPA Condensate Sample Port US Condensate
Sample Date Analysis/Sample I D		Conducted by	Level (MCL)	Level Source	5/19/2017 <b>20170705003</b>	5/8/2017 <b>20170705004</b>
Physical Characteristics	Units	Dy		Source	20170705005	20170705004
pH	pH units	U.S.	4.5-8.5	41000	7.43	7.98
Conductivity	μS/cm	U.S.			112	300
Anions (IC)						0.5
Fluoride Cations (IC)	mg/L	U.S.			0.6	0.5
Ammonia as Nitrogen (NH3-N)	mg/L	U.S.	1	SWEG& 41000	13.6	35.4
Trace Metals (ICP/MS)						
Calcium	mg/L	U.S.	30	41000	0.10	0.15
Sodium	mg/L	U.S. U.S.			0.30 17	0.34 10
Manganese	μg/L μg/L	U.S.	300	SWEG&41000	17	25
Nickel	μg/L	U.S.	300	SWEG& 41000	117	501
Silver	μg/L	U.S.	400	SWEG& 41000	12	18
Zinc	μg/L	U.S.	2,000	SWEG&41000	1,150	1,180
Silicon (ICP/MS) Silicon	μg/L	U.S.			3,720	10,500
Total Organic Carbon (Sievers)	µg/∟	0.0.			0,720	10,000
Inorganic Carbon	mg/L	U.S.			11.6	27.8
Organic Carbon	mg/L	U.S.	3	41000	41.4	48.1
Volatile Organics Acetone	μg/L	U.S.	15,000	SWEG	7,820	2,410
Volatile Organics - Special Interest Compounds (Semi-			15,000	SWEG	7,820	2,410
Trimethylsilanol	μg/L	U.S.			230	370
Semi-volatiles (GC/MS) - Target List						
Benzothiazole	μg/L	U.S.			56	59
N-n-Butylbenzenesulfonamide Tris(2-Chloroethyl)phosphate	μg/L μg/L	U.S. U.S.			60 <40	83 110
Decamethylcyclopentasiloxane	μg/L	U.S.			<40	102
Methyl sulfone	μg/L	U.S.	1,500,000	interim SWEG (06-2017)	58	163
Acid Extractables-EPA 625 List						
Benzoic acid Phenol	μg/L μg/L	U.S. U.S.	4,000	SWEG	<u>&lt;200</u> 179	389 <40
p-Cresol (4-Methylphenol)	μg/L μg/L	U.S.	4,000	SWEG	163	<40
Base/Neutral Extractables - EPA 625 List	P-3 -					
Benzyl alcohol	μg/L	U.S.			<40	1,490
Diethylphthalate	μg/L	U.S.			327	642
Semi-volatiles (GC/MS) - Special Interest Compounds 2-(2-Butoxyethoxy)ethanol	<b>(Semi-quar</b> μg/L	U.S.	Curve)		210	210
N,N-Dimethyl acetamide	μg/L μg/L	U.S.			240	550
N,N-Dimethylformamide	μg/L	U.S.			370	570
Dipropylene glycol methyl ether	μg/L	U.S.			140	330
2-Ethoxyethanol 2-Ethylhexanoic acid	μg/L	U.S. U.S.			<u>280</u> 110	230 not found
Ibuprofen	μg/L μg/L	U.S. U.S.			970	not found
1-Methyl-2-pyrrolidinone	μg/L	U.S.			<160	280
Monomethyl phthalate	μg/L	U.S.			110	110
(+)-Neomenthol	μg/L	U.S.			81	56
2-Phenoxyethanol 2-Phenyl-2-propanol	μg/L μg/L	U.S. U.S.			64 <80	55 160
1,3,5-Triallyl-1,3,5-triazine-2,4,6(1H,3H,5H)-trione	μg/L μg/L	U.S.			<80	80
Tributyl phosphate	μg/L	U.S.			36	48
Alcohols (DAI/GC/MS)						
Ethanol Methanol	μg/L	U.S. U.S.	40,000	SWEG	<u>33,600</u> 9,740	<400 6,860
2-Propanol (Isopropanol)	μg/L μg/L	U.S. U.S.	40,000	SVVEG	<u>9,740</u> 518	6,860 <400
Glycols (DAI/GC/MS)	<u>~~</u>					
1,2-Ethanediol (Ethylene glycol)	μg/L	U.S.	4000	SWEG	2,020	1,950
1,2-Propanediol (Propylene glycol)	µg/L	U.S.	1,700,000	SWEG	5,820	14,100
Silanols (LC/RI) (R & D Method -NIST traceable sta Dimethylsilanediol (DMSD)	n <b>dard not a</b> v μg/L	vailable) U.S.	35,000	SWEG	12,000	37,000
Carboxylates (IC)	μy/L	0.3.	30,000	GVVEG	12,000	37,000
Acetate	μg/L	U.S.			530	<500
Organic Carbon Recovery	percent	U.S.			85.82	51.65
Unaccounted Organic Carbon	mg/L	U.S.			5.87	23.26

NA=Not analyzed MI=Matrix Interference N/A=Not applicable