

# TOBSWMF's Leachate Monitoring Program December 2020

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Town of Babylon Department of Environmental  
Control

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**March 2021**

Laboratory data and summary report from December 2020 sampling for Babylon's Leachate Monitoring Program.

## **TOBSWMF's Leachate Monitoring Program**

**December 2020**

As part of its solid waste infrastructure the Town of Babylon maintains four ashfills, the Southern Ashfill (SA), the Old Northern U Ashfill (ONU), the New Northern U Ashfill (NNU) and the lateral expansion of the Southern Ashfill, also known as Cell 7 (NYSDEC Permit No. 1-4720-00778/00014). These ash facilities are located on the northern and southern face of the former Babylon Landfill located on Glean Street in West Babylon, NY.

Babylon's leachate monitoring program (LMP) samples leachate from each of Babylon's ash facilities pursuant to the requirements of 6NYCRR part 363 (formerly part 360) and/or special condition attached to their NYSDEC solid waste management operating permits. Sampling procedures are described in detail within the 2018 Update Site Analytical Plan for the Town of Babylon Solid Waste Management Facilities (SAP) (TOBDEC, 2018).

Historically for the TOBSWMF's LMP, sampling at the SA, ONU and NNU ash facilities was limited to baseline parameters. In 2018 the NYSDEC required Babylon also sample for 1,4 dioxane and PFOA/PFAS when sampling these facilities for the LMP. December 2020 sampling for the LMP also included these emerging contaminants. Leachate at Cell 7 continues to be sampled for expanded parameters (the expanded parameters list was modified as part of the updated NYSDEC Solid Waste Management Facility regulations (appendix 2)). Sampling of the SA, ONU, NNU and Cell 7 were performed on December 10, 2020. The sampling protocol for the LMP is detailed in the Updated SAP for the Town of Babylon Solid Waste Management Facilities (TOBDEC, 2018). Sampling at the SA and ONU is limited to the Secondary Leachate Collection and Recovery System (SLCRS). Sampling at the NNU is performed for both the Primary Leachate Collection and Recovery System (PLCRS) and SLCRS. Sampling at Cell 7 was for the PLCRS. The complete laboratory report, case narrative and QA/QC package from Pace Analytical Services Inc has been attached as an appendix to this report. Included within the report prepared by Pace Labs is lab analysis for PFAS/PFOA's performed by Eurofins Environmental Testing America. In addition to internal laboratory QA/QC, a trip blank for VOC's was obtained as part of the operational QA/QC requirements. The trip blank was clean. The method blank provided as part of the PFAS/PFOA's analysis for Cell 7 was clean. The method blank provided with PFAS/PFOA analysis for the remaining leachate facilities was clean. An equipment blank and field duplicate were included in the Groundwater Monitoring Program performed concurrently with the LMP sampling. The results of the equipment blank and field duplicate were not notable.

Project narratives prepared by the laboratory for each category were reviewed. Notations and flagging qualifiers discussed in the narratives were noted. The narrative for baseline analysis

stated acetone at the NNUP facility could have a high bias. For the Cell 7 facility the narrative noted that sulfate was detected in the blank. Each data package was certified by the laboratory as being in compliance with the laboratories quality assurance manual both technically and for completeness.

This section of the LMP report provides a brief summary of the December 2020 leachate sampling at the TOBSWMF's. The sections that follow provide a more detailed discussion of the results from each ash facility.

The following are notable observations from the December 2020 LMP sampling results:

- Manganese (.56 mg/l) did not exceed its MCL at the ONU. Manganese has exceeded its MCL at the ONU in 22 of the past 35 sampling events.
- pH of leachate at the ONU was 8.02, 8.59 at the SA, 8.56 at the NNU PLCRS, 7.37 at the NNU SLCRS and 7.93 at Cell 7. All continue to be observed within an acceptable range.
- Baseline organics observed at each facility for the December 2020 LMP:
  - Carbon disulfide was the sole baseline organic observed at the ONU (.0033 mg/l) and SA (.0015 mg/l) facilities.
  - Total baseline organics observed at the NNU facility; 0.1794 mg/l at the NNU P and 0.6986 mg/l at the NNU S.
  - No individual organic compound from the baseline parameters list (SA, ONU and NNU), or summation of those compounds (TTO)<sup>1</sup> were observed at or above their MCL or TTO limits at any of these Babylon ash facilities during the December 2020 LMP.
- Total organics from the expanded parameters list (above mdl) observed at the Cell 7 facility was .063 mg/l. Total organics observed above their reporting limit was .054 mg/l. Total Toxic Organics (TTO) (>.01 mg/l) at the Cell 7 facility was .031 mg/l. This is below the overall TTO limit (10 mg/l) and 1.5 mg/l limit for acid extractable compounds within the Town of Babylon discharge Certificate issued by SCDPW.
- Barium did not exceed its MCL at the ONU, SA, NNU or Cell 7 for December 2020.
- Mercury was observed slightly above its reporting limit at the NNUP (.00022 mg/l). Mercury was not detected at the ONU, SA, NNU SLCRS or Cell 7 for December 2020.
- Piper diagrams for the SA, ONU, NNU and Cell 7 were updated with leachate sampled during the December 2020 LMP and conform to historical data.
- Project narratives were prepared by Pace Analytical Services Inc. for the December 2020 LMP laboratory results. Any issues, deficiencies or flagging of results were summarized in these narratives, and can be found in the appendix of this report. Each data package

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<sup>1</sup> Suffolk County Department of Public Works Total Toxic Organics (TTO) limited to: VOC's 2.5 mg/l, Base Neutral Extractable Compounds 1.5 mg/l, Acid Extractable Compounds 1.5 mg/l and Pesticides and PCB's 1 mg/l.

TOBDEC

was certified by the laboratory as being in compliance with its contract for Babylon's LMP both technically and for completeness.

## TOBSWMF's Leachate Monitoring Program

### Old Northern U

### December 2020

Pursuant to NYSDEC 6NYCRR Part 363 requirements for the operation of the Town of Babylon's Old Northern U (ONU) Ashfill, leachate from that facility's secondary leachate collection and recovery system (SLCRS) was sampled in accordance with the procedures detailed in the TOBSWMF's SAP (TOBDEC, 2018). The ONU SLCRS is sampled semi-annually for baseline parameters. Pursuant to NYSDEC requirement to sample for "emerging contaminants", Babylon expanded sampling to include 1,4 dioxane and PFAS/PFOA's for this facility beginning in December 2019.

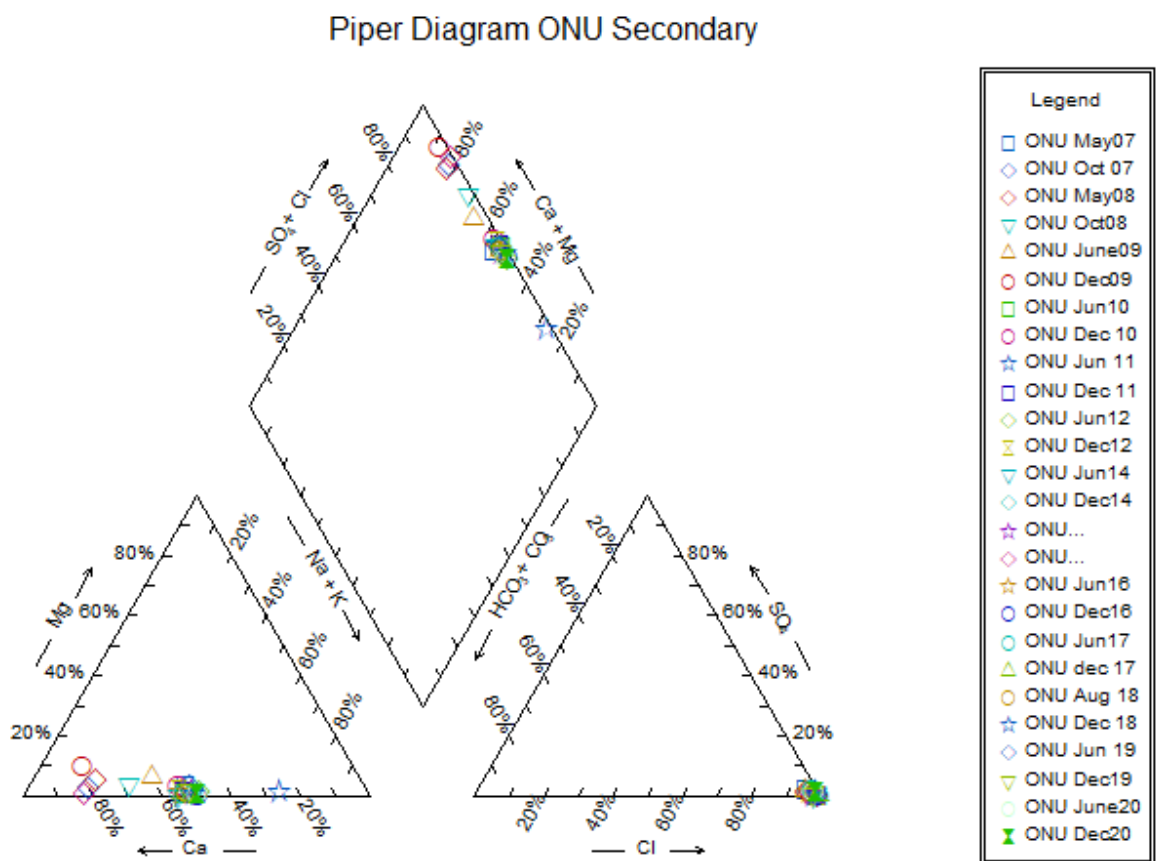
Ash has not been deposited in the ONU since it was capped in 2002 when the New Northern U (NNU) was constructed atop the facility. Leachate continues to be generated at the ONU despite the facility being capped and numerous attempts to locate the source. The LMP will continue at the ONU until there is a cessation of leachate generation. Included in this report is the December 2020 laboratory report from Pace Analytical Services, a spreadsheet summarizing parameters of concern dating back to 1995, a Piper diagram and a discussion of the laboratory results.

The attached spreadsheet provides a historical overview of leachate composition and any exceedance of MCL's at the ONU. The bullets below highlight notable observations from this round of sampling at the ONU and/or provide follow-up discussion/analysis of previous reports when appropriate.

- The chemical composition of leachate from the ONU for December 2020 generally conforms to historical data from the facility.
- pH measured in the field at the ONU SLCRS for June 2020 was 8.02.
- Manganese (.556 mg/l) was observed below its MCL for December 2020. Manganese has been observed exceeding its MCL in 22 of the past 35 monitoring events at the ONU.
- Barium (1.11 mg/l) was not observed above its MCL at the ONU for December 2020.
- Arsenic and lead were not detected above their mdl at the ONU for December 2020. Low values of arsenic and lead have been intermittently observed at this facility.
- Other metals observed at the ONU at values above their reporting limit and below their MCL (where one has been established) for December 2020 include boron (.385 mg/l), calcium (3360 mg/l), iron (1.55 mg/l), magnesium (7.27 mg/l), potassium (1360 mg/l), and sodium (3160 mg/l).

- Organic compounds from the baseline list observed at the ONU for December 2020 included Carbon disulfide (.0033 mg/l).
- 1,4 dioxane was observed at .57 ug/l for December 2020 at the ONU.
- Sulfide was not detected above its mdl at the ONU facility for December 2020.
- The Piper diagram from the ONU facility was updated with December 2020 data. The geochemical fingerprint for this facility remains unchanged.
- PFAS/PFOA’s results are attached in appendix 1.

The next round of sampling at the ONU is scheduled for June 2021.



Note: Solid hourglass = data point for December 2020.



PARAMETERS

03 MCL Oct\_08 June\_09 Dec\_09 June\_10 Dec\_10 Jun\_11 Dec\_11 12-Jun DEC\_12 Jun\_13 Dec\_13 Jun\_14 DEC\_14 June\_15 Dec\_15

perfluorobutanoic acid (PFBA)

perfluoropentanoic acid (PFPeA)

perfluorohexanoic acid(PFHxA)

perfluoroheptanoic acid

perfluorooctanoic acid(PFOA)

perfluorononanoic acid(PFNA)

perfluorodecanoic acid (PFDA)

perfluoroundecanoic acid(PFUnA)

perfluorododecanoic acid(PFDoA)

perfluorotridecanoic acid(PFTriA)

perfluorotetradecanoic acid(PFTeA)

perfluorobutanesulfonic acid(PFBS)

perfluorohexanesulfonic acid(PFHxS)

perfluoroheptanesulfonic acid(PFHpS)

perfluorooctanesulfonic acid(PFOS)

perfluorodecanesulfonic acid(PFDS)

perfluorooctanesulfonamide(FOSA)

N-methylperfluorooctanesulfonamidoacetic acid(NMeFOSAA)

N-ethylperfluorooctanesulfonamidoacetic acid(NEtFOSAA)

6:2FTS

8:2FTS



PARAMETERS	Jun_16	Dec_16	17-Jun	Dec_17	Aug_18	Dec_18	Jun_19	Dec_19	Jun_20	Dec_20
<b>CHLORIDE</b>	D 9630	D 44600	9970	348000	16400	19600	20400	D 14600	11600	12300
<b>SULFATE</b>	D 165	D 58	282	93.8	264	257	D 197	D 141	191	208
<b>Alkalinity</b>	D 271	182	143	148	293	139	245	302	196	137
<b>Na</b>	2390	8460	2500	6760	3720	3760	D 4560	D 3140	2230	3160
<b>K</b>	945	3870	1030	3310	1320	1570	D 1560	D 1140	937	1360
<b>Ca</b>	2960	9220	3100	8040	4290	4220	5140	D 3550	2390	3360
<b>Mg</b>	38.5	<10	19.4	0.293	19.2	11	192	71	12	7.27
<b>pH</b>	5.74	9.59/7	6.49	9.8	7.49	7.52	7.22	7.59	7.15	8.02
<b>TDS</b>	23900	52800	25200	69200	28600	24000	29900	19500	13700	20900
<b>PHENOL</b>										
<b>PHENOLS</b>	<.005	0.297	0.0264	0.0587	0.134	0.0059	<.00001	0.0158	<.005	0.0054
<b>IRON</b>	4.79	<5	4.32	<.4	2.21	1.44	31.8	13.3	6.16	1.55
<b>MANGANESE</b>	5.07	<.5	1.63	<.01	1.23	0.62	41.8	14.5	1.3	0.556
<b>TKN</b>	13.7	64.3	12.6	52.2	37.3	13.3	27.1	29.1	11.2	14
<b>ALUMINUM</b>	0.0704	J <10	<.0134	1.13	<10	<.2	<.2	<.2	<.2	<.2
<b>ACETONE</b>	J <	0.0804	<.001	0.0514	0.0024	J 0.0029	<.005	<.005	<.005	<.005
<b>3+4 methylphenol</b>										
<b>Methyl Ethyl Ketone</b>	<	<.005	<.0005	.0025	J <.005	<.005	<.005	<.005	<.005	<.005
<b>Arsenic</b>	<	<.5	<.0068	<.01	<.5	<.01	<.2	D <.01	<.01	<.01
<b>Lead</b>	0.0051	<.25	<.0013	<.4	<.25	0.0085	0.031	<.005	<.005	<.005
<b>Barium</b>	0.829	<10	1.32	4.9	1.34	J 1.13	2.77	2.07	0.619	1.11
<b>Xylene</b>	<	<.005	<.0005	<.002	<.003	<.003	<.003	<.003	<.003	<.003
<b>Zinc</b>	0.0358	<1	<.0012	<.02	<1	<.02	<.02	<.02	<.02	<.02
<b>Beryllium</b>	0.0022	J <.25	<.00057	.0036	J <.25	<.005	<.005	0.00034	J 0.00013	0.00017
<b>Nickel</b>	<	<2	<.00088	<.04	<2	<.04	<.04	<.04	0.0478	0.0193
<b>Selenium</b>	<	<.5	<.0062	<.01	<.5	<.01	<.2	D 0.0135	<.01	<.01
<b>Thallium</b>	<	<.5	<.0036	<.01	<.5	0.0085	J 0.0798	<.01	<.01	<.01
<b>Silver</b>	B <	<.5	<.0036	<.01	<.5	<.01	0.0048	J 0.0035	J 0.0047	<.01
<b>Toluene</b>	<	<.005	<.0005	<.001	<.001	<.001	<.001	<.001	<.001	<.001
<b>Carbon Disulfide</b>	<	<.005	<.0005	<.001	<.001	<.001	<.001	<.001	<.001	0.0033
<b>methylene chloride</b>	<	<.005	<.0005	<.001	<.001	<.01	<.001	<.001	<.001	<.001
<b>chromium</b>	<	<.5	<.0016	<.01	<.5	<.01	0.0071	J 0.0074	J 0.0489	0.0077
<b>Antimony</b>	<	<3	<.003	<.06	<3	<.06	0.06	<.06	<.06	<.06
<b>4-Methyl-2-pentanone</b>	J <	<.005	<.0005	<.005	<.005	<.005	<.005	<.005	<.005	<.005
<b>Sulfide</b>	<20	<2	<.61	9.6	<2	<.002	8	<2	1.6	<2
<b>1,4 dioxane</b>					0.21	JH 0.66	21	18.6	0.38	0.57

PARAMETERS	Jun_16	Dec_16	17-Jun Dec_17	Aug_18	Dec_18	Jun_19	Dec_19	Jun_20	Dec_20
perfluorobutanoic acid (PFBA)							180	B 73	76
perfluoropentanoic acid (PFPeA)							120	43	67
perfluorohexanoic acid(PFHxA)							160	60	82
perfluoroheptanoic acid							53	25	29
perfluorooctanoic acid(PFOA)							150	44	48
perfluorononanoic acid(PFNA)							17	7.3	8.1
perfluorodecanoic acid (PFDA)							5.4	J 2.1	1.8 J
perfluoroundecanoic acid(PFUnA)							ND	ND	nd
perfluorododecanoic acid(PFDoA)							ND	ND	nd
perfluorotridecanoic acid(PFTriA)							ND	ND	nd
perfluorotetradecanoic acid(PFTeA)							ND	ND	nd
perfluorobutanesulfonic acid(PFBS)							76	51	82
perfluorohexanesulfonic acid(PFHxS)							69	B 13	B 17
perfluoroheptanesulfonic acid(PFHpS)							2.8	J 0.42	J 0.47 J
perfluorooctanesulfonic acid(PFOS)							98	32	29
perfluorodecanesulfonic acid(PFDS)							ND	ND	nd
perfluorooctanesulfonamide(FOSA)							ND	0.76	JE nd
N-methylperfluorooctanesulfonamidoaceti							ND	ND	nd
N-ethylperfluorooctanesulfonamidoacetic ;							ND	ND	nd
6:2FTS							ND	ND	nd
8:2FTS							ND	ND	nd

## TOBSWMF's Leachate Monitoring Program

### Southern Ashfill

### December 2020

Pursuant to NYSDEC 6NYCRR Part 363 (formerly part 360) requirements for the operation of the Town of Babylon's Southern Ashfill (SA), leachate from that facility's Secondary Leachate Collection and Recovery System (SLCRS) was sampled in accordance with the procedures detailed in the TOBSWMF's SAP (TOBDEC, 2018). The SA facility requires semiannual sampling of leachate for baseline parameters from the facility's SLCRS. Pursuant to NYSDEC requirement to sample for "emerging contaminants", Babylon expanded sampling to include 1,4 dioxane and PFAS/PFOA's for this facility beginning in December 2019. This report includes the laboratory report from Pace Analytical Services, a Piper diagram, a spreadsheet summarizing parameters of concern dating back to 1994, and a discussion of the results.

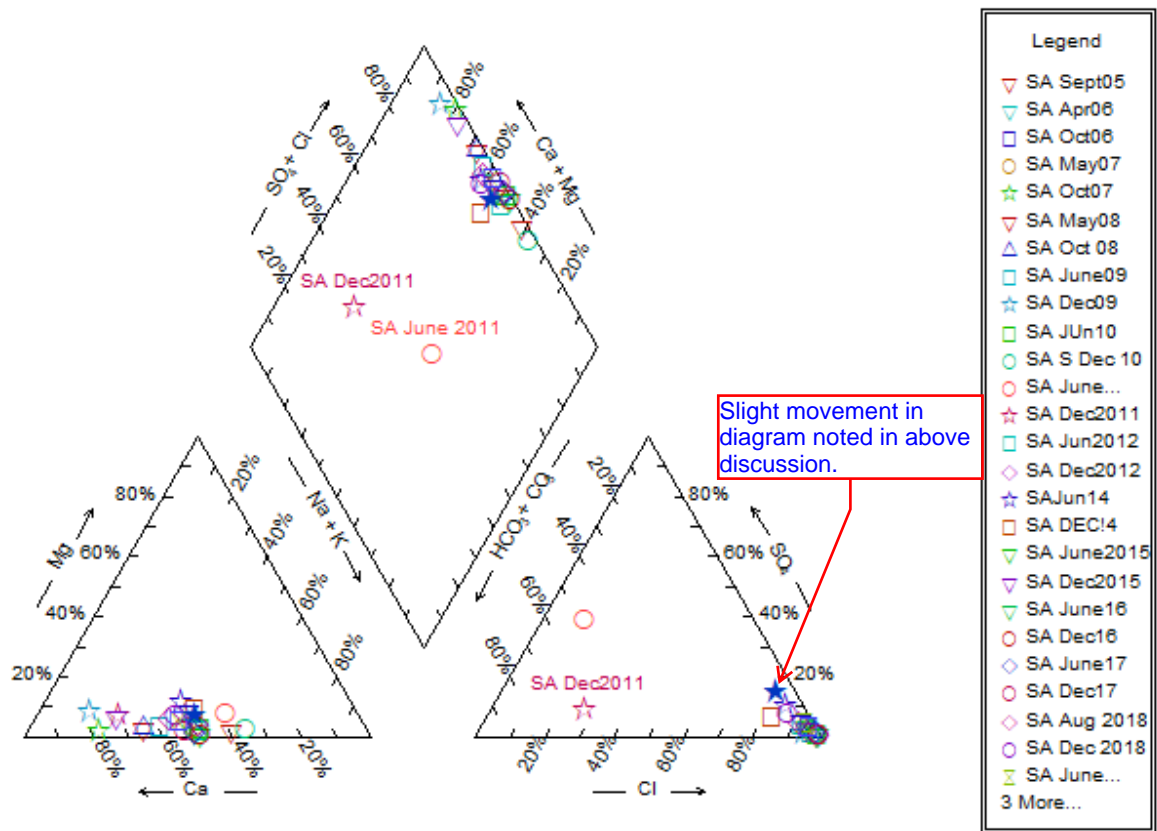
The attached spreadsheet provides a historical overview of leachate composition at the SA and any exceedance of the MCL's. The following bullets summarize any findings from this round of sampling at the SA and provide follow-up analysis or discussion when recommended from previous reports.

- Leachate indicators at the SA have been observed to be variable. Data from the December 2020 LMP at the SA fall within the range of historical data.
- A Piper diagram that includes SA data from December 2020 conforms to its established pattern. The slight movement in the lower right quadrant (highlighted on diagram) is likely the result of sulfate being observed slightly above its normal range and chloride being observed slightly below its normal range. These values while slightly outside their normal range have previously been observed during the facility's history.
- Lead (.098 mg/l) and arsenic (.015 mg/l) were observed at the SA for December 2020. Low values of lead and arsenic have been observed intermittently at the SA.
- Manganese was observed below its MCL at 4.21 mg/l for December 2020. Manganese had exceeded its MCL (8 mg/l) in June 2019. The only other sampling event where manganese exceeded its MCL at the SA facility was December 2013.
- Barium was observed at 0.189 mg/l at the SA for December 2020.
- Other metals observed at the SA at values above their reporting limit and below their MCL (where one has been established) for December 2020 include aluminum (3.86 mg/l), boron (.424mg/l), chromium (.0195 mg/l), calcium (642 mg/l), cadmium (.0036 mg/l), copper (.188 mg/l), iron (64.3 mg/l), magnesium (55.2 mg/l), nickel (.042 mg/l), potassium (247 mg/l), sodium (565 mg/l), and zinc (.486 mg/l).

- Carbon disulfide (.0015 mg/l) was the only organic from the baseline parameters list detected at the SA facility for December 2020.
- 1,4 dioxane was detected at 0.81 ug/l at the SA for December 2020.
- Mercury (.00011 mg/l) was detected below its reporting level at the SA for December 2020.
- pH measured in the field was 8.02 at the SA facility.
- Sulfide was not detected above its mdl at the SA facility for December 2020.
- PFAS/PFOA's results are attached in appendix A.

The next round of sampling is scheduled for June 2021.

Piper Diagram SA-Secondary LCRS



Note: Solid star indicates December 2020 data.











SA PARAMETERS	03 MCL	Dec_15	Jun_16	Dec_16	17-Jun	Dec_17	Aug_18	Dec_18	June_19	Dec_19	Jun_20	Dec_20
TKN	na	9.4700 D	3.8800	43.2000	28.4000	24.2000	0.5800	1.8000	17.0000 D	2.9	1.2	1.3
TDS	na	16600.0000	12.6000	39900	43000.0000	33200.0000	6130.0000	6300.0000	9360.0000	6800	8290	5250
Phenols	na	<.005	<.005	0.277	0.0124	0.0103	0.0569 J	0.0028 J	<.01	0.0092	<.005	0.0051
Chloride	na	6990.0000 D	#####	31100.0000	15400	57900.0000	3630.0000	2330	5830 D	5470	6860	2540
Sulfide	na		<20	<2	<.61	<2	<2	<2	6.4	<2	<2	<2
Iron	na	17.8000	2.3500	<5	6.86	11.7000	0.4540	12.8	210	2.85	21.5	64.3
Manganese	8 mg/l	4.97	1.87	<.5	3.42	3.86	2.09	1.09	8.44	5.31	6.67	4.21
Phenol	1.5 mg/l											
Xylene	2.5 mg/l *		<	<.005	<.0005	<.002	<.003	<.003	<.003	<.003	<.003	<.003
1,2,4 Trimethylbenzene	na											
SULFATE	na	263.0000 D	182.0000 D	246	221.0000	423.0000	251.0000	267.0000 D	361.0000 D	427	322	621
Arsenic	.4 mg/l	0.0048 B	<.01	<.5	<.0068	<.01	<.01	<.01	0.0599	<.01	<.01	0.0154
Acetone	na ppm	0.002 J	<	0.048	0.0755	0.0264	0.0032 J	<.005	0.0016 J	<.005	<.005	<.005
pH	5 - 12.5	7.0100	6.5300	7.21/6.5	6.18	6.95	8.08	8.05	8	7.24	8.12	8.59
Aluminum	na	0.0527 B	<	<10	<.0134	.0823 J	0.0506 J	0.564	13.5	<.2	0.531	3.86
Barium	8 mg/l	0.6040	0.4350	<10	1.62	1.08	0.205	0.17 J	0.481	0.158 J	0.264	0.189
Lead		0.0042	0.0023 J	<.25	<.0013	0.0058	0.0028 J	0.013	0.279	<.005	0.011	0.0982
Zinc		0.0109 B	0.1060	<1	0.0352	.0163 J	0.0097 J	0.0652	1.87	0.0064 J	0.0762	0.486
Toluene	2.5 mg/l *		<	<.005	<.0005	<.001	<.001	<.001	<.001	<.001	<.001	<.001
Cadmium	.8 mg/l	0.0011 B	<	<.125	<.000063	<.0025	<.0025	<.0025	0.0125	<.0025	<.0025	0.0036
Vanadium		<	<	<2.5	<.0008	<.05	<.05	0.0016 J	0.0226 J	<.05	<.05	0.0113 J
Tin												
Antimony		<	<	<3	<.003	<.06	<.06	<.06	0.0765	<.06	<.06	0.0252 J
Copper	1.6 mg/l	0.0073 B	0.0026 J	<1.25	<.0025	.011 J	0.0042 J	0.0185 J	0.36	0.0087 J	0.0374	0.188
Selenium	.4 mg/l	0.0026 B	<	<.5	<.0062	<.01	<.01	<.01	<.01	<.01	<.01	<.01
Silver	.4 mg/l	0.0035 B	<	<.5	<.0036	<.01	<.01	<.01	0.0043 J	0.0038 J	0.0028	<.01
Beryllium		<	0.0009 J	<.25	0.0051	.0018 J	<.005	<.05	<.005	0.00022 J	0.00011	<.005
Chromium	8 mg/l	0.0016 B	0.0414	<.5	<.0016	<.01	0.003 J	0.0067	0.0989	0.0156	0.0342	0.0195
Nickel	8 mg/l	0.0054 B	0.0243 J	<2	<.00088	<.04	<.04	<.04	0.069	<.04	0.0352	0.0415
Thallium		0.0244	<	<.5	<.0036	.0025 J	<.01	<.01	0.0276	0.012	<.01	<.01
Carbon disulfide			<	<.005	<.0005	<.001	<.001	<.001	<.001	<.001	<.001	0.0015
Methylene Chloride	2.5 mg/l		<	<.005	<.0005	<.001	<.001	<.001	<.001	<.001	<.001	<.001
Alkalinity		261 D	178	151	206	149	225	223	183	268	199	244
Ammonia		1.28	4.39 D	57.1000	11.8	26.9	0.05 J	0.75	4.7	2.9	0.23	0.00097 J
Hardness		4700 D	3400 D	16400.0000	11800	9600	2500	2200	4000	4000	10000	1400
1,4 dioxane	ug/l						0.37 J	0.75	0.88	<.2	0.9	0.81

SA PARAMETERS	03 MCL	Dec_15	Jun_16	Dec_16	17-Jun Dec_17	Aug_18	Dec_18	June_19	Dec_19	Jun_20	Dec_20	
		Dec_15										
Chloride												
Sulfate												
Alkalinity												
Na		329	1170	1494.3944	4180	3360	560	538	1330	843	1200	565
K		640	520	1087.2889	1770	1750	305	293	486	324	418	247
Ca		1820	1410	2053.8333	4660	4420	914	807	1760	991	1550	642
Mg		99.5	63.1	70.7500	70.6	83.7	56.4	64.2	103	105	90.6	55.2
pH		7.01	7.01		6.18	6.95	8.08	8.05	8	7.24	8.12	8.59
perfluorobutanoic acid (PFBA)									70	B	76	37
perfluoropentanoic acid (PFPeA)									110		82	48
perfluorohexanoic acid(PFHxA)									130		130	58
perfluoroheptanoic acid									52		44	30
perfluorooctanoic acid(PFOA)									130		110	86
perfluorononanoic acid(PFNA)									11		11	9.8
perfluorodecanoic acid (PFDA)									15		19	18
perfluoroundecanoic acid(PFUnA)									ND	ND		nd
perfluorododecanoic acid(PFDoA)									ND		0.95 J	nd
perfluorotridecanoic acid(PFTriA)									ND	ND		nd
perfluorotetradecanoic acid(PFTeA)									ND	ND		nd
perfluorobutanesulfonic acid(PFBS)									23		36	15
perfluorohexanesulfonic acid(PFHxS)									36	B	46 B	14
perfluoroheptanesulfonic acid(PFHpS)									ND		2.8	1.2 J
perfluorooctanesulfonic acid(PFOS)									51		110	57
perfluorodecanesulfonic acid(PFDS)									ND	ND		nd
perfluorooctanesulfonamide(FOSA)									ND		0.38 JB	nd
N-methylperfluorooctanesulfonamidoacetic acid(NMeFOSAA)									ND	ND		nd
N-ethylperfluorooctanesulfonamidoacetic acid(NEtFOSAA)									ND	ND		nd
6:2FTS									6.3	J	11 J	nd
8:2FTS									ND	ND		0.73 J

## **TOBSWMF's Leachate Monitoring Program**

### **New Northern U Ashfill**

**December 2020**

Pursuant to NYSDEC 6NYCRR Part 363 (formerly part 360) requirements for the operation of the Town of Babylon's New Northern U Ashfill (NNU), leachate from the NNU Primary and Secondary Leachate Collection and Recovery System (PLCRS and SLCRS) were sampled in accordance with the procedures detailed in the TOBSWMF's Updated SAP (TOBDEC, 2018). These facilities are sampled semi-annually for baseline parameters as part of Babylon's Leachate Monitoring Program (LMP). Pursuant to NYSDEC requirement to sample for "emerging contaminants", Babylon expanded sampling to include 1,4 dioxane and PFAS/PFOA's for this facility beginning in December 2019. This document includes the laboratory report from Pace Analytical Services, Inc., a spreadsheet summarizing parameters of concern at the facility, a Piper diagram of leachate from each liner system, and a discussion of the results.

The NNU which began accepting ash in 2003 sits atop the ONU, separated by a double liner system, with each layer consisting of a bentonite blanket, liner and geocomposite. The NNU SLCRS is also separated from the ONU by the ONU cap. Both systems serve as near impermeable barriers. The elevation of the NNU system (approximately 25-30 feet above the water table) prevents groundwater infiltration from being considered a source of leachate to the system.

The attached spreadsheet provides a historical overview of leachate composition at the NNU, highlighting any exceedance of an MCL from the facility's PLCRS and SLCRS. The following discussion summarizes any noteworthy findings from the December 2020 sampling and provides follow-up analysis or discussion wherever necessary or recommended in previous reports.

- For the December 2020 LMP pH was 7.37 at the NNU SLCRS and 8.56 at the NNU PLCRS.
- The overall leachate characteristics of the NNU PLCRS and SLCRS largely conform to the historical dataset for this facility.
- Arsenic and lead were not observed above their reporting limit at the NNU SLCRS and NNU PLCRS for December 2020. Low values of arsenic and lead have been intermittently observed at this facility.
- Mercury was not observed above its mdl at the NNU PLCRS or NNU SLCRS for December 2020.
- Organics from the baseline parameters list observed at the NNU for December 2020 were limited to low concentrations of acetone, MEK, 4 methyl 2 pentanone and carbon

disulfide. Acetone was observed at .173 mg/l at the NNU PLCRS and .617 mg/l at the NNU SLCRS. Low concentrations of acetone have been observed at this facility since June 2010.

MEK was detected at the NNU PLCRS at .0064 mg/l and .0718 mg/l at NNU-SLCRS during December 2020 sampling. Trace values of MEK have been intermittently observed at this facility.

4 methyl 2pentanone was observed at .0056 mg/l at the NNUS. Trace values of 4 methyl 2 pentanone have been observed intermittently at the NNU facility.

Carbon disulfide was observed at .0034 mg/l at the NNUS. Carbon disulfide has been observed intermittently at this facility.

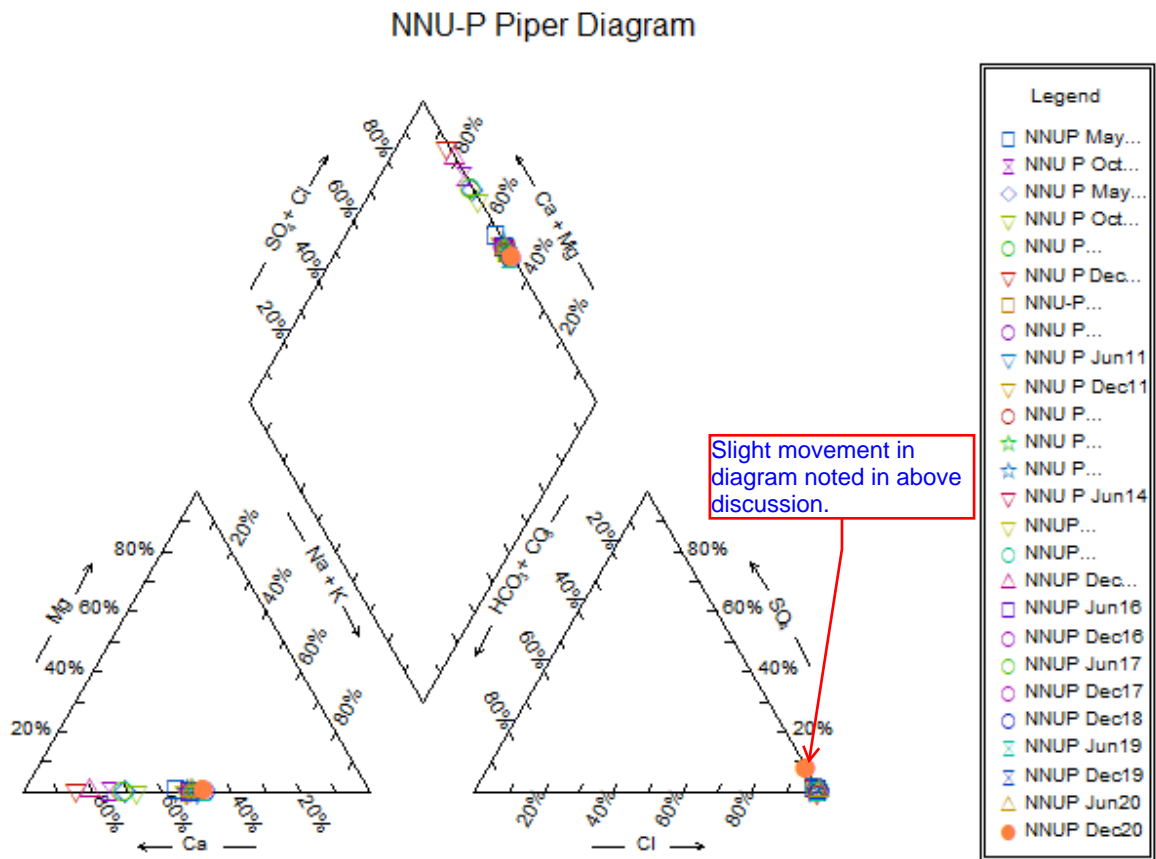
TTO as defined on the Town of Babylon discharge certificate issued by Suffolk County Department of Public Works is <.01 mg/l at the NNU facility.

Total baseline organics for the NNU PLCRS was .1794 mg/l and .6986 mg/l at the NNU SLCRS.

- 1,4 dioxane was observed at 2.3 ug/l at the NNU PLCRS and 2.6 ug/l at the NNU SLCRS.
- Barium was observed below its MCL at the NNU PLCRS (1.59 mg/l) and NNU SLCRS (2.42 mg/l) for December 2020. Barium has been observed exceeding its MCL at the NNU PLCRS 5 times over 35 sampling events through the life of the facility. Barium has exceeded its MCL at the NNU SLCRS 3 times over 35 sampling events through the life of the facility. The last exceedance for barium at each of the facilities was December 2012.
- Other metals observed above their reporting limit and below their MCL at the NNU PLCRS for December 2020 include Aluminum (.236 mg/l) boron (4.34 mg/l), cadmium (.0194 mg/l), chromium (.03 mg/l), calcium (6400 mg/l), copper (.211 mg/l), iron (.247 mg/l), magnesium (16 mg/l), manganese (.53 mg/l), nickel (.04 mg/l), potassium (2780 mg/l), sodium (6420 mg/l) and zinc (.096 mg/l).
- Other metals observed above their mdl and below their MCL at the NNU SLCRS for December 2020 include boron (5.63 mg/l), calcium (11100 mg/l), chromium (.02 mg/l), iron (.185mg/l), magnesium (2.3 mg/l), manganese (.312 mg/l), potassium (4680 mg/l), and sodium (10400 mg/l).
- Sulfide was below its mdl at both the NNUP and NNUS for the December 2020 LMP. Sulfide has exceeded its MCL at the NNUP for six of ten sampling rounds since June 2016. At the NNUS sulfide has also exceeded its MCL for 6 of 10 sampling rounds.
- BOD was observed below its MCL (300 mg/l) at the NNUP and NNUS. BOD has intermittently exceeded its MCL at these facilities.

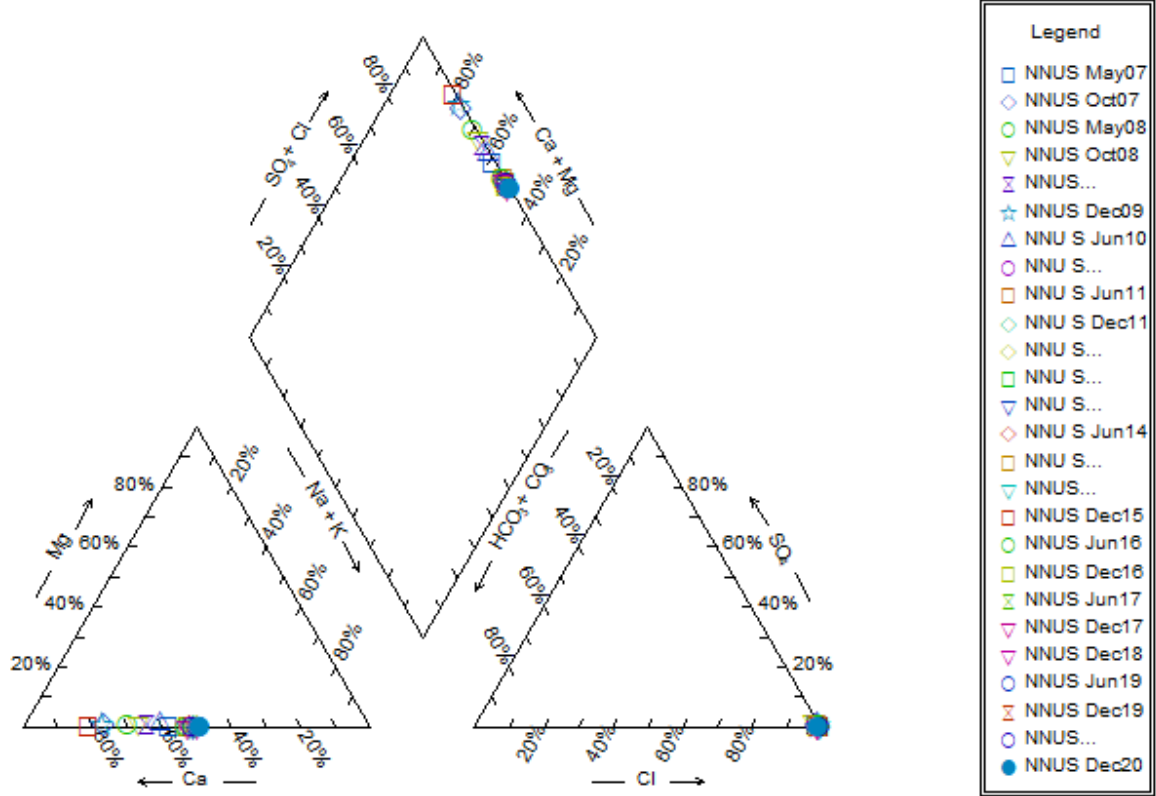
- A Piper diagram was prepared with the December 2020 data added to the historical dataset. The overall geochemical fingerprint for the NNU facilities is unchanged. A slight movement to the lower right quadrant of the NNUP facility is noted with the December 2020 data. This is likely the result of sulfate at the NNUP being observed above its normal range (this is not the highest value of sulfate observed at this facility).
- PFAS/PFOA's results are attached in appendix A.

The next round of sampling is scheduled for June 2021.



Note: solid circle represents December 2020 data.

Piper Diagram-NNU Secondary



Note: solid circle represents December 2020 data.



NNUP PARAMETERS

95 MCL

03 MCL

Aug\_03

Mar\_04

Sept\_04

Mar\_05

Sept\_05

Apr\_06

Oct\_06

May\_07

Oct\_07

May\_08

Oct\_08

June\_09

Dec\_09

Jun\_10

DEC\_10

Jun\_2011

DEC\_11

June\_12

perfluorodecanoic acid (PFDA)

perfluoroundecanoic acid(PFUnA)

perfluorododecanoic acid(PFDoA)

perfluorotridecanoic acid(PFTriA)

perfluorotetradecanoic acid(PFTeA)

perfluorobutanesulfonic acid(PFBS)

perfluorohexanesulfonic acid(PFHxS)

perfluoroheptanesulfonic acid(PFHpS)

perfluorooctanesulfonic acid(PFOS)

perfluorodecanesulfonic acid(PFDS)

perfluorooctanesulfonamide(FOSA)

N-methylperfluorooctanesulfonamidoacetic acit(NMeFOSAA)

N-ethylperfluorooctanesulfonamidoacetic acit(NEtFOSAA)

6:2FTS

8:2FTS







NNUP PARAMETERS	Jun_20	Dec_20
<b>CHLORIDE</b>	60800	24500
<b>SULFATE</b>	24.7	<b>2670</b>
<b>Alkalinity</b>	172	80.3
<b>Na</b>	10100	6420
<b>K</b>	4040	2780
<b>Ca</b>	11300	6400
<b>Mg</b>	2	16
<b>pH</b>	6.96	8.56
<b>hardness</b>		16200
<b>TDS</b>	70300	39900
<b>PHENOL</b>		
<b>PHENOLS</b>	0.118	0.0302
<b>IRON</b>	0.368	0.247
<b>MANGANESE</b>	0.335	0.53
<b>TKN</b>	107	61.1
<b>ALUMINUM</b>	<1	0.236
<b>ACETONE</b>	0.375	0.173
<b>Methyl Ethyl Ketone</b>	0.0412	0.0064
<b>Arsenic</b>	<.05	<.01
<b>Lead</b>	0.368	0.0047 J
<b>Barium</b>	2.24	1.59
<b>Cadmium</b>	<.0125	0.0194
<b>Copper</b>	0.103	0.211
<b>Selenium</b>	<.052	<.01
<b>Zinc</b>	0.12	0.096
<b>Carbon disulfide</b>	<.001	<.001
<b>BOD</b>	184	89.7
<b>Antimony</b>	<.3	0.026 J
<b>Beryllium</b>	0.00085	<.005
<b>Chromium</b>	0.237	0.0301
<b>Nickel</b>	0.111	0.0401
<b>Thallium</b>	0.0528	<.01
<b>Vanadium</b>	<.25	<.05
<b>methylene chloride</b>	<.001	<.001
<b>Toluene</b>	<.001	<.001
<b>Mercury</b>	<.0002	<.0002
<b>4-Methyl-2-pentanone</b>	0.0052	<.005
<b>Iodomethane</b>	0.0043	<.004
<b>sulfide mg/l</b>	<b>17.6</b>	<2
<b>1,4 Dioxane</b>	2.9	2.3
perfluorobutanoic acid (PFBA)	270	150
perfluoropentanoic acid (PFPeA)	130	120
perfluorohexanoic acid(PFHxA)	190	170
perfluoroheptanoic acid	31	39
perfluorooctanoic acid(PFOA)	43	54
perfluorononanoic acid(PFNA)	2.5	4.5

NNUP PARAMETERS	Jun_20	Dec_20
perfluorodecanoic acid (PFDA)	0.66 J	1.8 J
perfluoroundecanoic acid(PFUnA)	ND	nd
perfluorododecanoic acid(PFDoA)	ND	nd
perfluorotridecanoic acid(PFTriA)	ND	nd
perfluorotetradecanoic acid(PFTeA)	Nd	nd
perfluorobutanesulfonic acid(PFBS)	230	190
perfluorohexanesulfonic acid(PFHxS)	14 B	9.9
perfluoroheptanesulfonic acid(PFHpS)	0.29 J	.19 J
perfluorooctanesulfonic acid(PFOS)	12	12
perfluorodecanesulfonic acid(PFDS)	ND	nd
perfluorooctanesulfonamide(FOSA)	0.7 JB	nd
N-methylperfluorooctanesulfonamidoacetic acid	ND	nd
N-ethylperfluorooctanesulfonamidoacetic acid	ND	nd
6:2FTS	4.3 J	6
8:2FTS	ND	.59 J











NNUSPARAMETERS	95 MCL	Jun_20	Dec_20
<b>CHLORIDE</b>	500mg/l	61600	45900
<b>SULFATE</b>	500mg/l	8.9	65.9 J
<b>Alkalinity</b>		140	176
<b>Na</b>		10600	10400
<b>K</b>		4300	4680
<b>Ca</b>		11900	11100
<b>Mg</b>		2.02	2.3
<b>pH</b>	6.5-8.5	6.66	7.37
<b>TDS</b>	1000 mg/l	70800	71200
<b>PHENOL</b>	0.002mg/l		
<b>PHENOLS</b>		0.104	0.256
<b>IRON</b>	0.6mg/l	0.108	0.185
<b>MANGANESE</b>	0.6mg/l	0.322	0.312
<b>TKN</b>	10 mg/l	106	113
<b>ALUMINUM</b>	2mg/l	<1	0.0446 J
<b>ACETONE</b>	5 ppb	0.333	0.617
<b>Methyl Ethyl Ketone</b>	5 ppb	0.0406	0.0718
<b>Arsenic</b>	50 ppb	<.05	<.01
<b>Lead</b>	50 ppb	<.025	<.005
<b>Barium</b>		2.3	2.42
<b>Cadmium</b>		<.0125	<.0025
<b>Copper</b>		0.0895	<.025
<b>Zinc</b>		0.0405	<.02
<b>Antimony</b>		<.3	<.06
<b>Beryllium</b>		0.00089	<.005
<b>Chromium</b>		0.232	0.0201
<b>Nickel</b>		0.112	0.0281 J
<b>Selenium</b>		0.0468	<.01
<b>Thallium</b>		0.054	<.01
<b>Vanadium</b>		<.25	0.0062 J
<b>Silver</b>		0.0287	<.01
<b>methylene chloride</b>		<.001	<.001
<b>ammonia</b>		98.3	107
<b>hardness</b>		30800	32000
<b>carbon disulfide</b>		0.0018	0.0034
<b>4methyl2pentano</b>	ppb	0.0056	0.0064
<b>2 hexanone</b>		<.005	<.005
<b>Iodomethane</b>		0.0043	<.004
<b>sulfide</b>	12 mg/l	<b>12.8</b>	<2
<b>BOD</b>	300 mg/l	180	167
<b>1,4 dioxane</b>	ug/l	2.7	2.6
perfluorobutanoic acid (PFBA)		270	210
perfluoropentanoic acid (PFPeA)		130	150
perfluorohexanoic acid(PFHxA)		190	170
perfluoroheptanoic acid		30	29
perfluorooctanoic acid(PFOA)		36	28

NNUSPARAMETERS	95 MCL	Jun_20	Dec_20
perfluorononanoic acid(PFNA)		1.8 J	1.6 J
perfluorodecanoic acid (PFDA)		0.72 J	.58 J
perfluoroundecanoic acid(PFUnA)		ND	nd
perfluorododecanoic acid(PFDoA)		ND	nd
perfluorotridecanoic acid(PFTriA)		ND	nd
perfluorotetradecanoic acid(PFTeA)		ND	nd
perfluorobutanesulfonic acid(PFBS)		240	280
perfluorohexanesulfonic acid(PFHxS)		12 B	12
perfluoroheptanesulfonic acid(PFHpS)		ND	nd
perfluorooctanesulfonic acid(PFOS)		9.1	6.9
perfluorodecanesulfonic acid(PFDS)		ND	nd
perfluorooctanesulfonamide(FOSA)		3.1 B	nd
N-methylperfluorooctanesulfonamidoacetic acit(NMeFOSAA)		ND	nd
N-ethylperfluorooctanesulfonamidoacetic acit(NEtFOSAA)		ND	nd
6:2FTS		3.5 J	2.7 J
8:2FTS		ND	nd

## TOBSWMF's Leachate Monitoring Program

### Cell 7

#### December 2020

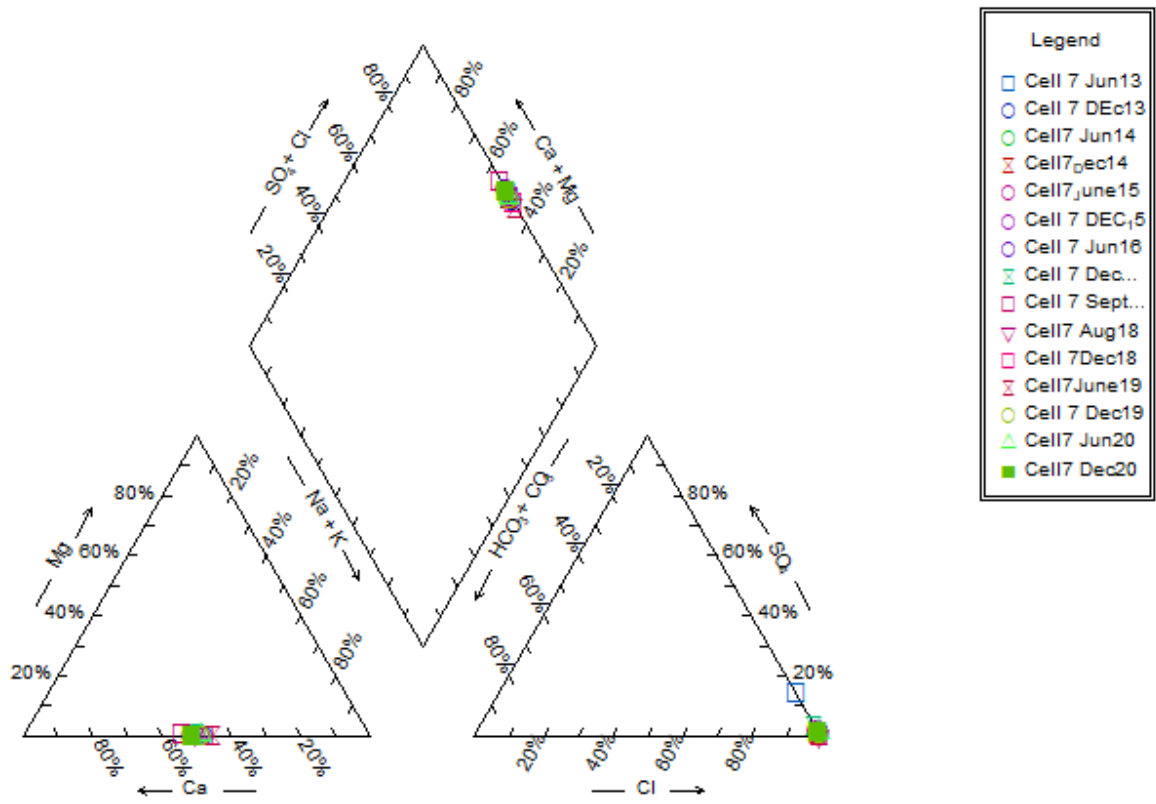
Pursuant to the NYSDEC operating permit for the operation of the Cell 7 Ashfill (Cell 7), leachate from that facility's PLCRS was sampled in accordance with the procedures detailed in the TOBSWMF's SAP (TOBDEC, 2018). The Cell 7 operating permit requires semiannual sampling of leachate for expanded parameters from the facility's PLCRS. The expanded parameters list is found within 6NYCRR part 363-4.6(h) and includes 1,4 dioxane, fluorinated alkyl substances (PFOA's) and other parameters (appendix 2) not found previously in NYCRR part 360. This report includes the laboratory report from Pace Analytical Services Inc., a spreadsheet summarizing the results, a Piper diagram and brief discussion.

- The overall leachate characteristics of the Cell 7 facility largely conform to the historical dataset for this facility.
- A Piper diagram that includes the December 2020 data for the Cell 7 facility was prepared and is attached to this section. The geochemical fingerprint as illustrated on the Piper diagram is unchanged for Cell 7.
- For December 2020 pH at Cell 7 was measured at 7.93.
- Analysis for 2378 TCDD and 2378 TCDF from December 2020 was ND (Reporting limit 10 pg/l).
- Analysis for 1,4 dioxane for December 2020 was reported at 1.7 ug/l.
- Mercury was not detected above its mdl at Cell 7 for December 2020.
- Organics from the expanded parameters list observed during December 2020 included acetone (.0497 mg/l), benzyl alcohol (.0031 mg/l), MEK (.0036mg/l), phenol (.0031 mg/l), 2,4 D (.0014 mg/l), 3-4 methylphenol (.0012 mg/l), carbon disulfide (.0011 mg/l), 2,4,5 T (.0013 mg/l) and endrin aldehyde (.00003 mg/l). Total expanded organics observed Cell 7 above their mdl for December 2020 was .063 mg/l.
- TTO (>.01 mg/l) observed at the Cell 7 facility for December 2020 is .0031 mg/l (phenol). This is below the overall TTO limit of 10 mg/l, and below the limit for acid extractable organic compounds of 1.5 mg/l set forth in the Town of Babylon Discharge Certification issued by SCDPW.
- No metals were observed above their MCL. Metals observed above their RL include aluminum (.311 mg/l), barium (3.16 mg/l), boron (.594 mg/l), calcium (6120 mg/l), chromium (.011 mg/l), iron (.702 mg/l), magnesium (7.17 mg/l), manganese (.255 mg/l), potassium (3160 mg/l) and sodium (4860 mg/l).

- PFAS/PFOA and 1,4 dioxane results are included in appendix 1.

The next round of sampling for leachate at the Cell 7 facility is scheduled for June 2021.

Piper Diagram Cell 7 PLCRS



Note: solid green square represents December 2020 data.

Cell7 PLCRS

CELL 7 PLCRS														
				07/01/13	3/13/2014	3/13/2014	06/25/14	12/12/14	06/16/15	12/14/2015				
				7/1/2013	13-Dec	DUP_1213	6/25/2014	12/12/2014	6/16/2015	12/14/2015	6/20/2016	Jan-17	Sept_17	Dec_17
TestNo	Analyte	CAS	Units											
	pH				7.88	1/30/2014	5.91	6.93	6.95		6.01	8.21	6.48	
	DO		mg/l		2.24	1/30/2014	1.31	0.86	1.77		0.87	1.87	0.53	
	Spec cond				61484		50900	45794	48822		56196	25443	65674	
	ORP						-256.4	-281.9	-276.2		-79.5	11.5	-326.5	
SW8270C	Pyrene	129-00-0	µg/L	10 U	10 U		ND U	ND U	ND U		10U	<2.5		<5.0
SW8270C	Safrole	94-59-7	µg/L	10 U	10 U		ND U	ND U	ND U	10 U	10U	<2.5		<5.0
SW9014	Cyanide	57-12-5	UG/L	10.0 U	10 U		50.0 U	10 U	20 U	10 U	10U	<2.9	<10	
SW9060	Total Organic Carbon		mg/L	51.6 D	108 D		35.2	88.0 D	21.3	2.5	22.6	<0.63	43.2	
E1613	Dioxin		Pg/L	1.0 U	10 U		ND	ND	ND U	10 U	10 U			
E300.0	Bromide	24959-67-9	mg/L	308 D	336 D			311 D	ND U	230 D	248D	117	373	
E300.0	Sulfate	14808-79-8	mg/L	5140 D	55 D		157 D	270 D	720 D	364 D	329D	338	375	
E351.2	Nitrogen, Kjeldahl, Total		mg/L	63.6 D	95 D		85.0 D	61.2 D	49.7 D	52.0 D	57.2D	17.1	67	
E353.2	Nitrate as N	14797-55-8	mg/L	2.50 U	2.00 U		2.00 U	0.100 U	0.100 U	0.10 U	.1U	<0.0050	<.05	
E353.2	Nitrite as N	14797-65-0	mg/L	0.100 U	0.100 U		0.100 U	0.100 U	0.100 U	0.10 U	.1U	<0.0050	<.05	
E410.4	Chemical Oxygen Demand		mg/L	517 D	1220 D		445 D	852 D	550 D	175 D	1400 D	560	1560	
E420.1	Phenolics, Total Recoverable		µg/L	49.4 D	309 D		66.6	47.5	54.8 D	5.0 U	41.9	76.2	110	
M3500-Cr D	Chromium, Hexavalent	18540-29-9	mg/L	0.0200 U	0.0200 U		0.0200 U	0.0200 U	0.0200 U	0.02 U	0.0200 U	<0.0030	<.1	
SM2120B	Color		units	75 D	150 D		200 D	150 D	75.0 D	15.0	25.0	40.0	25	
SM2320B	Alkalinity, Total (As CaCO3)		mg/L	181 D	266 D		223 D	273 D	175 D	119 D	122	78.6	160	
SM2340C	Hardness (As CaCO3)		mg/L	17200 D	13100 D		14200 D	17700 D	17800 D	13200 D	25800 D	6400	19600	
SM2540C	Total Dissolved Solids		mg/L	93900 D	39300 D		49400	51700	74000	55500	61100	2960	74800	
SM4500-CL	Chloride	16887-00-6	mg/L	23500 D	21600 D		21800 D	27900 D	26500 D	18400 D	18600 D	8320	31600	
SM4500-NH	Nitrogen, Ammonia (As N)	7664-41-7	mg/L	55.8 D	89.5 D		79.0 D	58.1 D	63.9 D	46.3 D	66.5 D	16.3	56.4	
SM5210B	Biochemical Oxygen Demand		mg/L	42	101		30	266	25	10 U	4	<3.3	43.5	
SW6010B	Aluminum	7429-90-5	UG/L	190 U	28.0 B		43.9 B	200 U	17.6 BN	39.5 B	200 U	200 U		
SW6010B	Antimony	7440-36-0	UG/L	24.0 U	4.0 B		15.8 B	60.0 U	13.2 BN	10.9 B	15.7 J	20.3 J		
SW6010B	Arsenic	7440-38-2	UG/L	56.0 U	8.4 B		39.0	19.1	11.4 N	21.1	19.9	7.6 J		
SW6010B	Barium	7440-39-3	UG/L	3170 B	2430		3490	2750	3940	2790	4250	954		
SW6010B	Beryllium	7440-41-7	UG/L	2.0 U	0.14 U		0.091 U	5.00 U	0.15 U	0.20 U	1.4 J	0.61 J		
SW6010B	Boron	7440-42-8	UG/L	958 B	381		333	666	673	480	651	429		
SW6010B	Cadmium	7440-43-9	UG/L	2.0 U	0.11 U		0.14 U	5.00 U	0.16 U	0.10 U	2.5 U	2.8	<2.5	
SW6010B	Calcium	7440-70-2	UG/L	6610000	6300000		7460000	7100000 D	7360000	5490000 DE	8830000	2570000	7180000	
SW6010B	Chromium	7440-47-3	UG/L	8.0 U	3.2 B		3.8 B	10.0 U	2.8 B	41.9	10 U	10 U		
SW6010B	Cobalt	7440-48-4	UG/L	8.0 U	0.19 U		0.16 U	50.0 U	1.5 B	0.20 U	50 U	2.6 J		
SW6010B	Copper	7440-50-8	UG/L	90.0 B	13.1 B		4.3 B	28.9	0.37 U	4.0 B	10.4 J	25 U		
SW6010B	Iron	7439-89-6	UG/L	896 B	839		1560	1480	894	3110	1230	1680	260	
SW6010B	Lead	7439-92-1	UG/L	20.0 U	10.6		7.7	3.00 U	0.85 UN	1.3 UN	5.8	<50	<100	
SW6010B	Magnesium	7439-95-4	UG/L	9900 B	3710 B		4560 B	7160	8620	9510	10400	8040	24000	
SW6010B	Manganese	7439-96-5	UG/L	2640	1690		2300	852	2100	672	755	304	861	
SW6010B	Nickel	7440-02-0	UG/L	6.0 U	0.34 U		0.29 U	40.0 U	2.8 B	0.30 U	40 U	3.1 J		
SW6010B	Potassium	7440-09-7	UG/L	2990000	3570000		3910000	3990000 D	3860000	2900000 D	4170000	1270000	415000	
SW6010B	Selenium	7782-49-2	UG/L	46.0 U	2.2 B		1.7 B	5.00 U	2.7 UN	2.2 UN	10 U	10 U		
SW6010B	Silver	7440-22-4	UG/L	4.0 U	0.43 U		0.37 U	10.0 U	0.87 UN	0.50 U	10 U			
SW6010B	Sodium	7440-23-5	UG/L	6310000	5760000		6490000	6240000 D	6230000	4870000 DE	7100000	2190000	6730000	
SW6010B	Thallium	7440-28-0	UG/L	38.0 U	1.3 U		4.6 B	10.0 U	1.0 U	1.9 U	10 U	10 U		
SW6010B	Tin	7440-31-5	UG/L	14.0 U	3.7 B		7.7 B	40.0 U	6.6	3.4 B	3.2 J	50 U		
SW6010B	Vanadium	7440-62-2	UG/L	6.0 U	6.4 B		3.7 B	50.0 U	5.4 B	5.0 B	50 U	1.6 J		

Cell7 PLCRS

CELL 7 PLCRS														
				07/01/13	3/13/2014	3/13/2014	06/25/14	12/12/14	06/16/15	12/14/2015				
				7/1/2013	13-Dec	DUP_1213	6/25/2014	12/12/2014	6/16/2015	12/14/2015	6/20/2016	Jan-17	Sept_17	Dec_17
SW6010B	Zinc	7440-66-6	UG/L	6.0 U	8.7 B		11.5 B	154	12.8 BN	1.6 U	4.2 J	20 U		
SW7470	Mercury	7439-97-6	UG/L	0.18 B	1.2 B		0.10 U	0.3	0.10 U	0.10 U	0.20 U	<0.2	.039J	
SW8081/808	4,4'-DDD	72-54-8	µg/L	ND U	ND U		ND U	ND U	0.10 U	0.10 U	0.10 U	.1 U		<0.10
SW8081/808	4,4'-DDE	72-55-9	µg/L	ND U	ND U		ND U	ND U	0.10 U	0.10 U	0.10 U	.1 U		<0.10
SW8081/808	4,4'-DDT	50-29-3	µg/L	ND U	ND U		ND U	ND U	0.10 U	0.10 U	0.10 U	.1 U		<0.10
SW8081/808	Aldrin	309-00-2	µg/L	ND U	ND U		ND U	ND U	0.050 U	0.050 U	0.050 U	.05 U		<0.050
SW8081/808	alpha-BHC	319-84-6	µg/L	ND U	ND U		ND U	ND U	0.050 U	0.050 U	0.050 U	.05 U		<0.050
SW8081/808	Aroclor 1016	12674-11-2	µg/L	ND U	ND U		ND U	ND U	1.0 U	1.0 U	1.0 U	1 U		<1.0
SW8081/808	Aroclor 1221	11104-28-2	µg/L	ND U	ND U		ND U	ND U	2.0 U	2.0 U	2.0 U	2 U		<2.0
SW8081/808	Aroclor 1232	11141-16-5	µg/L	ND U	ND U		ND U	ND U	1.0 U	1.0 U	1.0 U	1 U		<1.0
SW8081/808	Aroclor 1242	53469-21-9	µg/L	ND U	ND U		ND U	ND U	1.0 U	1.0 U	1.0 U	1 U		<1.0
SW8081/808	Aroclor 1248	12672-29-6	µg/L	ND U	ND U		ND U	ND U	1.0 U	1.0 U	1.0 U	1 U		<1.0
SW8081/808	Aroclor 1254	11097-69-1	µg/L	ND U	ND U		ND U	ND U	1.0 U	1.0 U	1.0 U	1 U		<1.0
SW8081/808	Aroclor 1260	11096-82-5	µg/L	ND U	ND U		ND U	ND U	1.0 U	1.0 U	1.0 U	1 U		<1.0
SW8081/808	beta-BHC	319-85-7	µg/L	ND U	ND U		ND U	ND U	0.050 U	0.050 U	0.050 U	.05 U		0.14
SW8081/808	Chlordane	57-74-9	µg/L	ND U	ND U		ND U	ND U	1.0 U	1.0 U	1.0 U			
SW8081/808	delta-BHC	319-86-8	µg/L	ND U	ND U		ND U	ND U	0.050 U	0.050 U	0.050 U	.05 U		<0.050
SW8081/808	Dieldrin	60-57-1	µg/L	ND U	ND U		ND U	ND U	0.10 U	0.10 U	0.10 U	.1 U		<0.10
SW8081/808	Endosulfan I	959-98-8	µg/L	ND U	ND U		ND U	ND U	0.050 U	0.050 U	0.050 U	.05 U		<0.050
SW8081/808	Endosulfan II	33213-65-9	µg/L	ND U	ND U		ND U	ND U	0.10 U	0.10 U	0.10 U	.1 U		<0.10
SW8081/808	Endosulfan sulfate	1031-07-8	µg/L	ND U	ND U		ND U	ND U	0.10 U	0.10 U	0.10 U	.1 U		<0.10
SW8081/808	Endrin	72-20-8	µg/L	ND U	ND U		ND U	ND U	0.10 U	0.10 U	0.10 U	.1 U		<0.10
SW8081/808	Endrin aldehyde	7421-93-4	µg/L	ND U	ND U		ND U	ND U	0.10 U	0.10 U	0.10 U	.1 U		<0.10
SW8081/808	gamma-BHC	58-89-9	µg/L	ND U	ND U		ND U	ND U	0.050 U	0.050 U	0.050 U	.05 U		<0.050
SW8081/808	Heptachlor	76-44-8	µg/L	ND U	ND U		ND U	ND U	0.050 U	0.050 U	0.050 U	.05 U		0.61
SW8081/808	Heptachlor epoxide	1024-57-3	µg/L	ND U	ND U		ND U	ND U	0.050 U	0.050 U	0.050 U	.05 U		<0.050
SW8081/808	Methoxychlor	72-43-5	µg/L	ND U	ND U		ND U	ND U	0.50 U	0.50 U	0.50 U	.5 U		<0.50
SW8081/808	Toxaphene	8001-35-2	µg/L	ND U	ND U		ND U	ND U	5.0 U	5.0 U	5.0 U	5 U		<5.0
SW8141A	Dimethoate	60-51-5	µg/L	ND U	ND U		ND U	ND U	1.0 U	1.0 U	1.0 U	.96 U		<.96
SW8141A	Disulfoton	298-04-4	µg/L	ND U	ND U		ND U	ND U	1.0 U	1.0 U	1.0 U	.96 U		<.96
SW8141A	Methyl parathion	298-00-0	µg/L	ND U	ND U		ND U	ND U	1.0 U	1.0 U	1.0 U	.96 U		<.96
SW8141A	Parathion	56-38-2	µg/L	ND U	ND U		ND U	ND U	1.0 U	1.0 U	1.0 U	.96 U		<.96
SW8141A	Phorate	298-02-2	µg/L	ND U	ND U		ND U	ND U	1.0 U	1.0 U	1.0 U	.96 U		<.96
SW8141A	Thionazin	297-97-2	µg/L	ND U	10 U		ND U					<2.5		<5.0
SW8151	2,4,5-T	93-76-5	µg/L	ND U	ND U		ND U	0.25 U	0.25 U	0.25 U	0.25 U	.047 J		<0.25
SW8151	2,4,5-TP (Silvex)	93-72-1	µg/L	ND U	ND U		0.33 P	0.25 U	0.25 U	0.25 U	0.25 U	.25 U		<0.25
SW8151	2,4-D	94-75-7	µg/L	3.2 P	ND U		0.26 PJ	0.50 U	0.57 P	0.52 P	0.50 U	.5 U		0.28 J
SW8151	Dinoseb	88-85-7	µg/L	ND	ND U		ND U	1.3	0.37 P	0.76 P	0.20 U	.085 J		<0.20
SW8260B	1,1,1,2-Tetrachloroethane	630-20-6	µg/L	ND U	ND U		ND U	ND U	5.0 U	5 U	5.0 U	<0.50	<1.0	<1.0
SW8260B	1,1,1-Trichloroethane	71-55-6	µg/L	ND U	ND U	ND U	ND U	ND U	5.0 U	5 U	5.0 U	<0.50	<1.0	<1.0
SW8260B	1,1,2,2-Tetrachloroethane	79-34-5	µg/L	ND U	ND U	ND U	ND U	ND U	5.0 U	5 U	5.0 U	<0.50	<1.0	<1.0
SW8260B	1,1,2-Trichloroethane	79-00-5	µg/L	ND U	ND U	ND U	ND U	ND U	5.0 U	5 U	5.0 U	<0.50	<1.0	<1.0
SW8260B	1,1-Dichloroethane	75-34-3	µg/L	ND U	ND U	ND U	ND U	ND U	5.0 U	5 U	5.0 U	<0.50	<1.0	<1.0
SW8260B	1,1-Dichloroethene	75-35-4	µg/L	ND U	ND U	ND U	ND U	ND U	5.0 U	5 U	5.0 U	<0.50	<1.0	<1.0
SW8260B	1,1-Dichloropropene	563-58-6	µg/L	ND U	ND U	ND U	ND U	ND U	5.0 U	5 U	5.0 U	<0.50	<1.0	<1.0
SW8260B	1,2,3-Trichloropropane	96-18-4	µg/L	ND U	ND U	ND U	ND U	ND U	5.0 U	5 U	5.0 U	<0.50	<1.0	<1.0
SW8260B	1,2-Dibromo-3-chloropropane	96-12-8	µg/L	ND U	ND U	ND U	ND U	ND U	5.0 U	5 U	5.0 U	<0.50	<1.0	<1.0
SW8260B	1,2-Dibromoethane	106-93-4	µg/L	ND U	ND U	ND U	ND U	ND U	5.0 U	5 U	5.0 U	<0.50	<1.0	<1.0
SW8260B	1,2-Dichlorobenzene	95-50-1	µg/L	ND U	ND U	ND U	ND U	ND U	5.0 U	5 U	5.0 U	<0.50	<1.0	<1.0

Cell7 PLCRS

CELL 7 PLCRS				07/01/13	3/13/2014	3/13/2014	06/25/14	12/12/14	06/16/15	12/14/2015				
				7/1/2013	13-Dec	DUP_1213	6/25/2014	12/12/2014	6/16/2015	12/14/2015	6/20/2016	Jan-17	Sept_17	Dec_17
SW8260B	1,2-Dichloroethane	107-06-2	µg/L	ND U	ND U	ND U	ND U	ND U	5.0 U	5 U	5.0 U	<0.50	<1.0	<1.0
SW8260B	1,2-Dichloropropane	78-87-5	µg/L	ND U	ND U	ND U	ND U	ND U	5.0 U	5 U	5.0 U	<0.50	<1.0	<1.0
SW8260B	1,3-Dichlorobenzene	541-73-1	µg/L	ND U	ND U	ND U	ND U	ND U	5.0 U	5 U	5.0 U	<0.50	<1.0	<1.0
SW8260B	1,3-Dichloropropane	142-28-9	µg/L	ND U	ND U	ND U	ND U	ND U	5.0 U	5 U	5.0 U	<0.50	<1.0	<1.0
SW8260B	1,4-Dichlorobenzene	106-46-7	µg/L	ND U	ND U	ND U	ND U	ND U	5.0 U	5 U	5.0 U	<0.50	<1.0	<1.0
	1,4-Dioxane (p-Dioxane)		ug/l											<100
SW8260B	2,2-Dichloropropane	594-20-7	µg/L	ND U	ND U	ND U	ND U	ND U	5.0 U	5 U	5.0 U	<0.50	<1.0	<1.0
SW8260B	2-Butanone	78-93-3	µg/L	17	41 Z	39 DZ	23	35	16	5 U	5.0 U	<0.50	15.3	9.2
SW8260B	2-Hexanone	591-78-6	µg/L	ND U	ND U	ND U	ND U	ND U	5.0 U	5 U	5.0 U	<0.50	<5.0	<5.0
SW8260B	4-Methyl-2-pentanone	108-10-1	µg/L	1 J	3 J	3 DJ	2 J	2 J	1 J	5 U	5.0 U	<0.50	<5.0	1.3 J
SW8260B	Acetone	67-64-1	µg/L	120	260 E	270 D	110	300 E	110	5 U	5.0 U	15.6	209	77.1
SW8260B	Acetonitrile	75-05-8	µg/L	ND U	28	25 D	35	100	49	40	5.0 U	<2.5	<5.0	<5.0
SW8260B	Acrolein	107-02-8	µg/L	ND U	ND U	ND U	ND U	ND U	5.0 U	5 U	5.0 U	<0.50	<1.0	<1.0
SW8260B	Acrylonitrile	107-13-1	µg/L	ND U	ND U	ND U	ND U	ND U	5.0 U	5 U	5.0 U	<0.50	<1.0	<1.0
SW8260B	Allyl Chloride	107-05-1	µg/L	ND U	ND U	ND U	ND U	ND U	5.0 U	5 U	5.0 U	<0.50	<1.0	<1.0
SW8260B	Benzene	71-43-2	µg/L	ND U	ND U	ND U	ND U	ND U	5.0 U	5 U	5.0 U	<0.50	<1.0	<1.0
SW8260B	Bromochloromethane	74-97-5	µg/L	ND U	ND U	ND U	ND U	ND U	5.0 U	5 U	5.0 U	<0.50	<1.0	<1.0
SW8260B	Bromodichloromethane	75-27-4	µg/L	ND U	ND U	ND U	ND U	ND U	5.0 U	5 U	5.0 U	<0.50	<1.0	<1.0
SW8260B	Bromoform	75-25-2	µg/L	ND U	ND U	ND U	ND U	ND U	5.0 U	5 U	5.0 U	<0.50	<1.0	<1.0
SW8260B	Bromomethane	74-83-9	µg/L	ND U	ND U	ND U	ND U	ND U	5.0 U	5 U	5.0 U	<0.50	<1.0	<1.0
SW8260B	Carbon disulfide	75-15-0	µg/L	ND U	ND U	ND U	ND U	ND U	5.0 U	5 U	5.0 U	<0.50	<1.0	<1.0
SW8260B	Carbon tetrachloride	56-23-5	µg/L	ND U	ND U	ND U	ND U	ND U	5.0 U	5 U	5.0 U	<0.50	<1.0	<1.0
SW8260B	Chlorobenzene	108-90-7	µg/L	ND U	ND U	ND U	ND U	ND U	5.0 U	5 U	5.0 U	<0.50	<1.0	<1.0
SW8260B	Chloroethane	75-00-3	µg/L	ND U	ND U	ND U	ND U	ND U	5.0 U	5 U	5.0 U	<0.50	<1.0	<1.0
SW8260B	Chloroform	67-66-3	µg/L	ND U	ND U	ND U	ND U	ND U	5.0 U	5 U	5.0 U	<0.50	<1.0	<1.0
SW8260B	Chloromethane	74-87-3	µg/L	ND U	ND U	ND U	ND U	ND U	5.0 U	5 U	5.0 U	<0.50	<1.0	<1.0
SW8260B	Chloroprene	126-99-8	µg/L	ND U	ND U	ND U	ND U	ND U	5.0 U	5 U	5.0 U	<0.50	<1.0	<1.0
SW8260B	cis-1,2-Dichloroethene	156-59-2	µg/L	ND U	ND U	ND U	ND U	ND U	5.0 U	5 U	5.0 U	<0.50	<1.0	<1.0
SW8260B	cis-1,3-Dichloropropene	10061-01-5	µg/L	ND U	ND U	ND U	ND U	ND U	5.0 U	5 U	5.0 U	<0.50	<1.0	<1.0
SW8260B	Dibromochloromethane	124-48-1	µg/L	ND U	ND U	ND U	ND U	ND U	5.0 U	5 U	5.0 U	<0.50	<1.0	<1.0
SW8260B	Dibromomethane	74-95-3	µg/L	ND U	ND U	ND U	ND U	ND U	5.0 U	5 U	5.0 U	<0.50	<1.0	<1.0
SW8260B	Dichlorodifluoromethane	75-71-8	µg/L	ND U	ND U	ND U	ND U	ND U	5.0 U	5 U	5.0 U	<0.50	<1.0	<1.0
SW8260B	Ethyl Methacrylate	97-63-2	µg/L	ND U	ND U	ND U	ND U	ND U	5.0 U	5 U	5.0 U	<0.50	<1.0	<1.0
SW8260B	Ethylbenzene	100-41-4	µg/L	ND U	ND U	ND U	ND U	ND U	5.0 U	5 U	5.0 U	<0.50	<1.0	<1.0
SW8260B	Iodomethane	74-88-4	µg/L	ND U	ND U	ND U	ND U	ND U	5.0 U	5 U	2 J	<0.50	<1.0	<1.0
SW8260B	Isobutyl alcohol	78-83-1	µg/L	ND U	ND U	ND U	14 J	ND U	25 U	25 U	25 U			
SW8260B	Methacrylonitrile	126-98-7	µg/L	ND U	ND U	ND U	ND U	ND U	5.0 U	5 U	5.0 U	<0.50	<1.0	<1.0
SW8260B	Methyl Methacrylate	80-62-6	µg/L	ND U	ND U	ND U	ND U	ND U	5.0 U	5 U	5.0 U	<0.50	<1.0	<1.0
SW8260B	Methylene chloride	75-09-2	µg/L	ND U	ND U	ND U	ND U	ND U	5.0 U	5 U	5.0 U	<0.50	<1.0	<1.0
SW8260B	Propionitrile	107-12-0	µg/L	ND U	ND U	ND U	ND U	ND U	5.0 U	5 U	5.0 U	<2.0	<4.0	<4.0
SW8260B	Silane, methoxytrimethyl-		ug/L	5 JN										
SW8260B	Silanol, trimethyl-		ug/L	19 JN				15 JN		13 JN				
SW8260B	Styrene	100-42-5	µg/L	ND U	ND U	ND U	ND U	ND U	5.0 U	5 U	5.0 U	<0.50	<1.0	<1.0
SW8260B	Tetrachloroethene	127-18-4	µg/L	ND U	ND U	ND U	ND U	ND U	5.0 U	5 U	5.0 U	<0.50	<1.0	<1.0
SW8260B	Toluene	108-88-3	µg/L	ND U	ND U	ND U	ND U	ND U	5.0 U	5 U	5.0 U	<0.50	<1.0	<1.0
SW8260B	trans-1,2-Dichloroethene	156-60-5	µg/L	ND U	ND U	ND U	ND U	ND U	5.0 U	5 U	5.0 U	<0.50	<1.0	<1.0
SW8260B	trans-1,3-Dichloropropene	10061-02-6	µg/L	ND U	ND U	ND U	ND U	ND U	5.0 U	5 U	5.0 U	<0.50	<1.0	<1.0
SW8260B	trans-1,4-Dichloro-2-butene	110-57-6	µg/L	ND U	ND U	ND U	ND U	ND U	5.0 U	5 U	5.0 U	<0.50	<1.0	<1.0
SW8260B	Trichloroethene	79-01-6	µg/L	ND U	ND U	ND U	ND U	ND U	5.0 U	5 U	5.0 U	<0.50	<1.0	<1.0

Cell7 PLCRS

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				07/01/13	3/13/2014	3/13/2014	06/25/14	12/12/14	06/16/15	12/14/2015				
				7/1/2013	13-Dec	DUP_1213	6/25/2014	12/12/2014	6/16/2015	12/14/2015	6/20/2016	Jan-17	Sept_17	Dec_17
SW8260B	Trichlorofluoromethane	75-69-4	µg/L	ND U	ND U	ND U	ND U	ND U	5.0 U	5 U	5.0 U	<0.50	<1.0	<1.0
SW8260B	Trimethylsilyl fluoride+Sulfur diox		ug/L	220 JN										
SW8260B	Vinyl acetate	108-05-4	µg/L	ND U	ND U	ND U	ND U	ND U	5.0 U	5 U	5.0 U	<0.50	<1.0	<1.0
SW8260B	Vinyl chloride	75-01-4	µg/L	ND U	ND U	ND U	ND U	ND U	5.0 U	5 U	5.0 U	<0.50	<1.0	<1.0
SW8260B	Xylene (total)	1330-20-7	µg/L	ND U	ND U	ND U	ND U	ND U	5.0 U	5 U	5.0 U	<0.50	<2.0	<2.0
SW8270C	1,2,4,5-Tetrachlorobenzene	95-94-3	µg/L	ND U	ND U	ND U	ND U	ND U	10 U	10 U	10 U	<2.5		<5.0
SW8270C	1,2,4-Trichlorobenzene	120-82-1	µg/L	ND U	10 U	40 U	ND U	ND U	10 U	10 U	10 U	<2.5		<5.0
SW8270C	1,2-Dichlorobenzene	95-50-1	µg/L	ND U	10 U	ND U	ND U	ND U	10 U	10 U	10 U	<2.5		<5.0
SW8270C	1,3,5-Trinitrobenzene	99-35-4	µg/L	ND U	ND U	40 U	ND U	ND U	10 U	10 U	10 U	<2.5		<5.0
SW8270C	1,3-Dichlorobenzene	541-73-1	µg/L	ND U	10 U	ND U	ND U	ND U	10 U	10 U	10 U	<2.5		<5.0
SW8270C	1,3-Dinitrobenzene	99-65-0	µg/L	ND U	ND U	ND U	ND U	ND U	10 U	10 U	10 U	<2.5		<5.0
SW8270C	1,4-Dichlorobenzene	106-46-7	µg/L	ND U	10 U	ND U	ND U	ND U	10 U	10 U	10 U	<2.5		<5.0
SW8270C	1,4-Naphthoquinone	130-15-4	µg/L	ND U	ND U	ND U	ND U	ND U	10 U	10 U	10 U	<2.5		<5.0
SW8270C	1-Naphthylamine	134-32-7	µg/L	ND U	10 U	ND U	ND U	ND U	10 U	10 U	10 U	<2.5		<5.0
SW8270C	2,2'-oxybis(1-chloropropane)	108-60-1	µg/L	ND U	ND U	40 U	ND U	ND U	10 U	10 U	10 U	<2.5		<5.0
SW8270C	2,3,4,6-Tetrachlorophenol	58-90-2	µg/L	ND U	ND U	ND U	ND U	ND U	10 U	10 U	10 U	<2.5		<5.0
SW8270C	2,4,5-Trichlorophenol	95-95-4	µg/L	25 U	25 U	ND U	ND U	ND U	25 U	25 U	25 U	<2.5		<5.0
SW8270C	2,4,6-Trichlorophenol	88-06-2	µg/L	10 U	ND U	ND U	ND U	ND U	10 U	10 U	10 U	<2.5		<5.0
SW8270C	2,4-Dichlorophenol	120-83-2	µg/L	ND U	10 U	ND U	ND U	ND U	10 U	10 U	10 U	<2.5		<5.0
SW8270C	2,4-Dimethylphenol	105-67-9	µg/L	10 U	10 U	40 U	ND U	ND U	10 U	10 U	10 U	<2.5		<5.0
SW8270C	2,4-Dinitrophenol	51-28-5	µg/L	ND U	25 U	ND U	ND U	ND U	25 U	25 U	25 U	<5.0		<10.0
SW8270C	2,4-Dinitrotoluene	121-14-2	µg/L	ND U	ND U	ND U	ND U	ND U	10 U	10 U	10 U	<2.5		<5.0
SW8270C	2,6-Dichlorophenol	87-65-0	µg/L	10 U	ND U	40 U	ND U	ND U	10 U	10 U	10 U	<2.5		<5.0
SW8270C	2,6-Dinitrotoluene	606-20-2	µg/L	10 U	10 U	ND U	ND U	ND U	10 U	10 U	10 U	<2.5		<5.0
SW8270C	2-Acetylaminofluorene	53-96-3	µg/L	ND U	ND U	ND U	ND U	ND U	20 U	20 U	20 U	<2.5		<5.0
SW8270C	2-Chloronaphthalene	91-58-7	µg/L	10 U	10 U	ND U	ND U	ND U	10 U	10 U	10 U	<2.5		<5.0
SW8270C	2-Chlorophenol	95-57-8	µg/L	10 U	ND U	ND U	ND U	ND U	10 U	10 U	10 U	<2.5		<5.0
SW8270C	2-Methylnaphthalene	91-57-6	µg/L	10 U	ND U	ND U	ND U	ND U	10 U	10 U	10 U	<0.17		<5.0
SW8270C	2-Methylphenol	95-48-7	µg/L	ND U	10 U	40 U	ND U	ND U	10 U	10 U	10 U	<2.5		<5.0
SW8270C	2-Naphthylamine	91-59-8	µg/L	10 U	ND U	ND U	ND U	ND U	10 U	10 U	10 U	<2.5		<5.0
SW8270C	2-Nitroaniline	88-74-4	µg/L	25 U	25 U	100 U	ND U	ND U	25 U	25 U	25 U	<2.5		<5.0
SW8270C	2-Nitrophenol	88-75-5	µg/L	10 U	10 U	40 U	ND U	ND U	10 U	10 U	10 U	<2.5		<5.0
SW8270C	3,3'-Dichlorobenzidine	91-94-1	µg/L	ND U	ND U	80 U	ND U	ND U	20 U	20 U	20 U	<2.5		<5.0
SW8270C	3,3'-Dimethylbenzidine	119-93-7	µg/L	10 U	10 U	40 U	ND U	ND U	10 U	10 U	10 U	<2.5		<5.0
SW8270C	3-Methylcholanthrene	56-49-5	µg/L	ND U	10 U	ND U	ND U	ND U	10 U	10 U	10 U	<2.5		<5.0
SW8270C	3-Methylphenol/4-Methylphenol	12-03-3	µg/L	9 J	150	170 D	ND U	9 J	41	10 U	10 U			16.8
SW8270C	3-Nitroaniline	99-09-2	µg/L	ND U	25 U	ND U	ND U	ND U	25 U	25 U	25 U	<2.5		<5.0
SW8270C	4,6-Dinitro-2-methylphenol	534-52-1	µg/L	ND U	ND U	ND U	ND U	ND U	25 U	25 U	25 U	<5.0		<10.0
SW8270C	4-Aminobiphenyl	92-67-1	µg/L	20 U	ND U	80 U	ND U	ND U	20 U	20 U	20 U	<2.5		<5.0
SW8270C	4-Bromophenyl-phenylether	101-55-3	µg/L	10 U	ND U	40 U	ND U	ND U	10 U	10 U	10 U	<2.5		<5.0
SW8270C	4-Chloro-3-methylphenol	59-50-7	µg/L	10 U	10 U	ND U	ND U	ND U	10 U	10 U	10 U	<2.5		<5.0
SW8270C	4-Chloroaniline	106-47-8	µg/L	10 U	ND U	ND U	ND U	ND U	10 U	10 U	10 U	<2.5		<5.0
SW8270C	4-Chlorophenyl-phenylether	7005-72-3	µg/L	10 U	10 U	40 U	ND U	ND U	10 U	10 U	10 U	<2.5		<5.0
SW8270C	4-Nitroaniline	100-01-6	µg/L	25 U	ND U	100 U	ND U	ND U	25 U	25 U	25 U	<2.5		<5.0
SW8270C	4-Nitrophenol	100-02-7	µg/L	25 U	ND U	100 U	ND U	ND U	25 U	25 U	25 U	<5.0		<10.0
SW8270C	5-Nitro-o-toluidine	99-55-8	µg/L	10 U	10 U	40 U	ND U	ND U	10 U	10 U	10 U	<2.5		<5.0
SW8270C	7,12-Dimethylbenz(a)anthracene	57-97-6	µg/L	10 U	ND U	ND U	ND U	ND U	10 U	10 U	10 U	<2.5		<5.0
SW8270C	Acenaphthene	83-32-9	µg/L	10 U	ND U	ND U	ND U	ND U	10 U	10 U	10 U	<0.22		<5.0
SW8270C	Acenaphthylene	208-96-8	µg/L	10 U	ND U	ND U	ND U	ND U	10 U	10 U	10 U	<0.21		<5.0



## Cell7 PLCRS

CELL 7 PLCRS														
				07/01/13	3/13/2014	3/13/2014	06/25/14	12/12/14	06/16/15	12/14/2015				
				7/1/2013	13-Dec	DUP_1213	6/25/2014	12/12/2014	6/16/2015	12/14/2015	6/20/2016	Jan-17	Sept_17	Dec_17
SW8270C	Acetophenone	98-86-2	µg/L	10 U	10 U	40 U	ND U	ND U	10 U	10 U	10 U	<2.5		1.2 J
SW8270C	Anthracene	120-12-7	µg/L	10 U	ND U	ND U	ND U	ND U	10 U	10 U	10 U	<2.5		0.61 J
SW8270C	Benzo(a)anthracene	56-55-3	µg/L	10 U	ND U	40 U	ND U	ND U	10 U	10 U	10 U	<2.5		<5.0
SW8270C	Benzo(a)pyrene	50-32-8	µg/L	10 U	10 U	40 U	ND U	ND U	10 U	10 U	10 U	<2.5		<5.0
SW8270C	Benzo(b)fluoranthene	205-99-2	µg/L	10 U	10 U	40 U	ND U	ND U	10 U	10 U	10 U	<2.5		<5.0
SW8270C	Benzo(g,h,i)perylene	191-24-2	µg/L	ND U	ND U	40 U	ND U	ND U	10 U	10 U	10 U	<2.5		<5.0
SW8270C	Benzo(k)fluoranthene	207-08-9	µg/L	ND U	10 U	40 U	ND U	ND U	10 U	10 U	10 U	<2.5		<5.0
SW8270C	Benzyl alcohol	100-51-6	µg/L	1	ND U	40 U	ND U	4 J	10 U	10 U	10 U	<2.5		<5.0
SW8270C	Bis(2-chloroethoxy)methane	111-91-1	µg/L	ND U	ND U	40 U	ND U	ND U	10 U	10 U	10 U	<2.5		<5.0
SW8270C	Bis(2-chloroethyl)ether	111-44-4	µg/L	ND U	10 U	40 U	ND U	ND U	10 U	10 U	10 U	<2.5		<5.0
SW8270C	Bis(2-ethylhexyl)phthalate	117-81-7	µg/L	10 U	ND U	ND U	ND U	ND U	10 U	10 U	10 U	<2.5		1.0 J
SW8270C	Butyl benzyl phthalate	85-68-7	µg/L	10 U	ND U	ND U	ND U	ND U	10 U	10 U	10 U	<2.5		<5.0
SW8270C	Chlorobenzilate	510-15-6	µg/L	ND U	ND U	ND U	ND U	ND U	10 U	10 U	10 U	<2.5		<5.0
SW8270C	Chrysene	218-01-9	µg/L	ND U	ND U	40 U	ND U	ND U	10 U	10 U	10 U	<2.5		<5.0
SW8270C	Diallate	2303-16-4	µg/L	ND U	10 U	ND U	ND U	ND U	10 U	10 U	10 U	<2.5		<5.0
SW8270C	Dibenzo(a,h)anthracene	53-70-3	µg/L	ND U	ND U	ND U	ND U	ND U	10 U	10 U	10 U	<2.5		<5.0
SW8270C	Dibenzofuran	132-64-9	µg/L	ND U	10 U	40 U	ND U	ND U	10 U	10 U	10 U	<2.5		<5.0
SW8270C	Diethylphthalate	84-66-2	µg/L	10 U	ND U	ND U	ND U	ND U	10 U	10 U	10 U	<2.5		0.15 J
SW8270C	Dimethylphthalate	131-11-3	µg/L	ND U	10 U	40 U	ND U	ND U	10 U	10 U	10 U	<2.5		<5.0
SW8270C	Di-n-butyl phthalate	84-74-2	µg/L	10 U	10 U	40 U	ND U	ND U	10 U	10 U	10 U	<2.5		<5.0
SW8270C	Di-n-octyl phthalate	117-84-0	µg/L	10 U	ND U	ND U	ND U	ND U	10 U	10 U	10 U	<2.5		<5.0
SW8270C	Ethyl methanesulfonate	62-50-0	µg/L	ND U	10 U	ND U	ND U	ND U	10 U	10 U	10 U	<2.5		<5.0
SW8270C	Famphur	52-85-7	µg/L	10 U	10 U	40 U	ND U	ND U	10 U	10 U	10 U	<5.0		<10.0
SW8270C	Fluoranthene	206-44-0	µg/L	ND U	10 U	40 U	ND U	ND U	10 U	10 U	10 U	<2.5		<5.0
SW8270C	Fluorene	86-73-7	µg/L	ND U	10 U	40 U	ND U	ND U	10 U	10 U	10 U	<0.17		<5.0
SW8270C	Hexachlorobenzene	118-74-1	µg/L	ND U	ND U	ND U	ND U	ND U	10 U	10 U	10 U	<2.5		<5.0
SW8270C	Hexachlorobutadiene	87-68-3	µg/L	10 U	ND U	40 U	ND U	ND U	10 U	10 U	10 U			<5
SW8270C	Hexachlorocyclopentadiene	77-47-4	µg/L	ND U	ND U	40 U	ND U	ND U	10 U	10 U	10 U	<2.5		<5.0
SW8270C	Hexachloroethane	67-72-1	µg/L	ND U	10 U	ND U	ND U	ND U	10 U	10 U	10 U	<2.5		<5.0
SW8270C	Hexachloropropene	1888-71-7	µg/L	ND U	10 U	ND U	ND U	ND U	10 U	10 U	10 U	<2.5		<5.0
SW8270C	Indeno(1,2,3-cd)pyrene	193-39-5	µg/L	10 U	ND U	40 U	ND U	ND U	10 U	10 U	10 U	<2.5		<5.0
SW8270C	Isodrin	465-73-6	µg/L	10 U	ND U	40 U	ND U	ND U	10 U	10 U	10 U	<2.5		<5.0
SW8270C	Isophorone	78-59-1	µg/L	10 U	ND U	ND U	ND U	ND U	10 U	10 U	10 U	<2.5		<5.0
SW8270C	Isosafrole	120-58-1	µg/L	ND U	ND U	40 U	ND U	ND U	10 U	10 U	10 U	<2.5		<5.0
SW8270C	Kepone	143-50-0	µg/L	ND U	10 U	40 U	ND U	ND U	10 U	10 U	10 U	<5.0		<10.0
SW8270C	Methapyrilene	91-80-5	µg/L	ND U	10 U	40 U	ND U	ND U	10 U	10 U	10 U	<2.5		<5.0
SW8270C	Methyl methanesulfonate	66-27-3	µg/L	ND U	10 U	40 U	ND U	ND U	10 U	10 U	10 U	<2.5		<5.0
SW8270C	Naphthalene	91-20-3	µg/L	10 U	ND U	ND U	ND U	ND U	10 U	10 U	10 U	<0.18		<5.0
SW8270C	Nitrobenzene	98-95-3	µg/L	10 U	10 U	40 U	ND U	ND U	10 U	10 U	10 U	<2.5		<5.0
SW8270C	N-Nitrosodiethylamine	55-18-5	µg/L	10 U	ND U	ND U	ND U	ND U	10 U	10 U	10 U	<2.5		<5.0
SW8270C	N-Nitrosodimethylamine	62-75-9	µg/L	ND U	10 U	40 U	ND U	ND U	10 U	10 U	10 U	<2.5		<5.0
SW8270C	N-Nitroso-di-n-butylamine	924-16-3	µg/L	ND U	ND U	40 U	ND U	ND U	10 U	10 U	10 U	<2.5		<5
SW8270C	N-Nitroso-di-n-propylamine	621-64-7	µg/L	ND U	10 U	40 U	ND U	ND U	10 U	10 U	10 U	<2.5		<5
SW8270C	N-Nitrosodiphenylamine	86-30-6	µg/L	10 U	ND U	40 U	ND U	ND U	10 U	10 U	10 U	<2.5		<5.0
SW8270C	N-Nitrosomethylethylamine	10595-95-6	µg/L	10 U	10 U	40 U	ND U	ND U	10 U	10 U	10 U	<2.5		<5.0
SW8270C	N-Nitrosopiperidine	100-75-4	µg/L	ND U	ND U	40 U	ND U	ND U	10 U	10 U	10 U	<2.5		<5.0
SW8270C	N-Nitrosopyrrolidine	930-55-2	µg/L	10 U	10 U	40 U	ND U	ND U	10 U	10 U	10 U	<2.5		<5.0
SW8270C	O,O,O-Triethylphosphorothioate	126-68-1	µg/L	ND U	10 U	ND U	ND U	ND U	10 U	10 U	10 U	<2.5		<5.0
SW8270C	o-Toluidine	95-53-4	µg/L	ND U	10 U	40 U	ND U	ND U	10 U	10 U	10 U	<2.5		<5.0

Cell7 PLCRS

CELL 7 PLCRS				07/01/13	3/13/2014	3/13/2014	06/25/14	12/12/14	06/16/15	12/14/2015				
				7/1/2013	13-Dec	DUP_1213	6/25/2014	12/12/2014	6/16/2015	12/14/2015	6/20/2016	Jan-17	Sept_17	Dec_17
SW8270C	p-Dimethylaminoazobenzene	60-11-7	µg/L	10 U	ND U	40 U	ND U	ND U	10 U	10 U	10 U	<2.5		<5.0
SW8270C	Pentachlorobenzene	608-93-5	µg/L	ND U	ND U	ND U	ND U	ND U	10 U	10 U	10 U	<2.5		<5.0
SW8270C	Pentachloronitrobenzene	82-68-8	µg/L	ND U	10 U	40 U	ND U	ND U	10 U	10 U	10 U	<2.5		<5.0
SW8270C	Pentachlorophenol	87-86-5	µg/L	ND U	25 U	100 U	ND U	ND U	25 U	25 U	25 U	<5.0		<10.0
SW8270C	Phenacetin	62-44-2	µg/L	10 U	ND U	ND U	ND U	ND U	10 U	10 U	10 U	<2.5		<5.0
SW8270C	Phenanthrene	85-01-8	µg/L	ND U	10 U	40 U	ND U	ND U	10 U	10 U	10 U	<0.17		<5.0
SW8270C	Phenol	108-95-2	µg/L	20	10 U	40 U	ND U	34	6 J	10 U	10 U	<2.5		19.4
SW8270C	p-Phenylenediamine	106-50-3	µg/L	10 U	10 U	ND U	ND U	ND U	10 U	10 U	10 U			<5.0
SW8270C	Pronamide	23950-58-5	µg/L	10 U	10 U	ND U	ND U	ND U	10 U	10 U	10 U	<2.5		<5.0
	Sulfide	18496-25-8	mg/L		2.00 U		2.00 U	25.3	2 U		20 U	<0.61	6.4	
EPA1613B	2378-TCDF		pg/l				ND		2 U					ND
EPA1613B	2378-TCDD		pg/l				ND		2 U		10 U			ND
ASTM D517	Total Uranium	7440-61-1	ng/l											1.07 ± 0.050 (0.193) C:NA T:NA
EPA 537	Perfluorobutanesulfonic acid PFBS	375-73-5	ng/l											<84
EPA 537	Perfluoroheptanoic acid PFHpA	375-85-9	ng/l											23
EPA 537	Perfluorohexanesulfonic acid PFHxS	355-46-4	ng/l											13 J
EPA 537	Perfluorononanoic acid PFNA	375-95-1	ng/l											<19
EPA 537	Perfluorooctanesulfonic acid PFOS	1763-23-1	ng/l											<38
EPA 537	Perfluorooctanoic acid PFOA	335-67-1	ng/l											29
EPA 903.1	Radium-226	13982-63-3	ng/l											3.02 ± 1.28 (1.13) C:NA T:33%
EPA 904.0	Radium-228	15262-20-1	ng/l											4.14 ± 1.79 (2.70) C:75% T:16%
	6:2 FTS		ng/l											
	8:2 FTS		ng/l											
	N-ethyl perfluorooctandsulfamidoacetic acidNEtFOSAA		ng/l											
	N-methylperfluorooctansulfamicacetic acid NMeFOSAA		ng/l											
	perfluorobutanoic acid PFBA		ng/l											
	perfluorodecansulfonic acid PFDS		ng/l											
	perfluorodecanoic acid PFDA		ng/l											
	perfluorododecanoic acid PFDoA		ng/l											
	perfluoroheptanesulfonic acid PFHps		ng/l											
	perfluorohexanoic acid PFHxA		ng/l											
	perfluorooctane sulfonamide FOSA		ng/l											
	perfluoropentanoic acid PFPeA		ng/l											
	perfluorotetradecanoic acid PFTeA		ng/l											
	perfluorotridecnaoic acid PFTriA		ng/l											
	perfluoroundecanoic acid PFUnA		ng/l											
	n-Nitrosomorpholine													
	Dimethylbenz(A) Anthracene													
	Bis(2-chloroisopropyl)ether													
	total PFOA/PFAS													

Cell7 PLCRS

CELL 7 PLCRS						
	Aug_18	Dec_18	Jun_19	Dec_19	June_20	Dec_20
Analyte						
pH	7.11	7.43	7.81	7.48	7.36	7.93
DO	0.05	2.01	0	1.7	2.59	2.02
Spec cond	788	1112	876	2194	>20,000	>20,000
ORP	-55.8	-75.1	-96.3	-79.2	-73.9	-102.4
Pyrene	U	<5	<5.0	<.25	<5	<5.0
Safrole	U	<5	<5.0	<.25	<5	<5.0
Cyanide	<10	21.3	4.6J	7	4.7	3.2J
Total Organic Carbon	94.7	84.8	257 D	147	69.2	28.8
Dioxin						
Bromide	353	350	516	422	480	260
Sulfate	10.3	6.5	7.2	335	129 D	305J D
Nitrogen, Kjeldahl, Total	51.2	56.3	104 D	65.2 D	93.8 D	21.6
Nitrate as N	<.05	0.051	0.090	<0.50 D	<0.050	<0.25 D
Nitrite as N	<.05	<.05	<0.050	<0.050	<0.050	<0.050
Chemical Oxygen Demand	1810	1690	3870	3410	2240	1120
Phenolics, Total Recoverable	236	177		358 D	278 D	35.3
Chromium, Hexavalent	<.1D	<.02	<.02	<.02	<.02	0.052
Color		15		50.0		250 D
Alkalinity, Total (As CaCO3)	275	216	336	223	176	123
Hardness (As CaCO3)	20400	20100	28800	26700	28400	15800
Total Dissolved Solids	54000	54400	74600	62000	58800	34000
Chloride	30500	29600	50600	48500	49500	22700
Nitrogen, Ammonia (As N)	51.7D	29.8	93.3	78.7	82.2	50.7 D
Biochemical Oxygen Demand	137D	134	494	235	103	46.8 D
Aluminum	<10000 D	<200	<1000 D	77.6J D	<1000 D	311
Antimony	<3000 D	18.8J	<300 D	45.4J D	<300 D	19.2J
Arsenic	<500 D	<10.0	<50.0 D	28.4 D	<50.0 D	<10.0
Barium	3580J D	3130	6450 D	5840 D	5550 D	3160
Beryllium	<250 D	<5.0	1.7J D	<10.0 D	0.58J D	0.20J
Boron	612J D	718	334 D	1040 D	92.5J D	594
Cadmium	<125 D	14.4J D	<12.5 D	<5.0 D	<12.5 D	<2.5
Calcium	8140000 D	7430000	9750000 D	9300000 D	9900000 D	6120000 D
Chromium	<500 D	<10.0	46.1J D	<20.0 D	157 D	11.4
Cobalt	<2500 D	5.0J	<250 D	<100 D	<250 D	<50.0
Copper	<1250 D	<25.0	59.0J D	<50.0 D	56.0J D	<25.0
Iron	10600 D	362	150 D	388 D	109 D	702
Lead	<250 D	<50.0 D	<25.0 D	<10.0 D	<25.0 D	<5.0
Magnesium	18100 D	11400	4420 D	11100 D	6450 D	7170
Manganese	3250 D	649	1440 D	750 D	221 D	255
Nickel	<2000 D	<40.0	<200 D	<80.0 D	72.0J D	26.2J
Potassium	3930000 D	4600000 D	6390000 D	5700000 D	5550000 D	3160000 D
Selenium	<500 D	<10.0	125 D	17.8J D	<50.0 D	<10.0
Silver	<500 D	<10.0	<50.0 D	<20.0 D	18.8J D	<10.0
Sodium	6910000 D	6870000 D	9900000 D	7950000 D	8800000 D	4860000 D
Thallium	<500 D	4.5J	<50.0 D	<20.0 D	<50.0 D	<10.0
Tin	<2500 D	<50.0	<250 D	<100 D	<250 D	<50.0
Vanadium	<2500 D	<50.0	<250 D	13.6J D	<250 D	10.0J

Cell7 PLCRS

CELL 7 PLCRS						
	Aug_18	Dec_18	Jun_19	Dec_19	June_20	Dec_20
Zinc	<1000 D	16.8J D	132 D	<40.0 D	<100 D	<20.0
Mercury	<.2	<0.20	0.15J	0.15J	<.2	<0.200
4,4'-DDD	<0.10	<0.10	<0.10	<0.10	<.1	<0.10
4,4'-DDE	<0.10	<0.10	<0.10	<0.10	<.1	<0.10
4,4'-DDT	<0.10	<0.10	<0.10	<0.10	<.1	0.023J
Aldrin	<0.050	<0.050	<0.050	<0.050	<.05	<0.050
alpha-BHC	<0.050	<.05	<0.050	<.05	<.05	<0.050
Aroclor 1016	<1.0	<1.0	<1.0	<1.0	<1	<1.0
Aroclor 1221	<2.0	<2.0	<2.0	<2.0	<1	<1.0
Aroclor 1232	<1.0	<1.0	<1.0	<1.0	<1	<1.0
Aroclor 1242	<1.0	<1.0	<1.0	<1.0	<1	<1.0
Aroclor 1248	<1.0	<1.0	<1.0	<1.0	<1	<1.0
Aroclor 1254	<1.0	<1.0	<1.0	0.68J	<1	<1.0
Aroclor 1260	<1.0	<1.0	<1.0	<1.0	<1	<1.0
beta-BHC	<.05	<.05	<0.050	<.05	<.05	<0.050
Chlordane						
delta-BHC	<.05	<.05	<0.050	<.05	<.05	<0.050
Dieldrin	<0.10	<0.10	<0.10	<0.10	<.1	<0.10
Endosulfan I	<0.050	<0.050	<0.050	<0.050	<.05	<0.050
Endosulfan II	<0.10	<0.10	<0.10	<0.10	<.1	<0.10
Endosulfan sulfate	<0.10	<0.10	<0.10	<0.10	<.1	<0.10
Endrin	<0.10	<0.10	<0.10	<0.10	<.1	<0.10
Endrin aldehyde	<0.10	<0.10	<0.10	<0.10	<.1	0.026J
gamma-BHC	<.05	<.05	<0.050	<.05	<.05	<0.050
Heptachlor	<.05	<0.050	<0.050	<0.050	<.05	<0.050
Heptachlor epoxide	<0.050	<0.050	<0.050	<0.050	<.05	<0.050
Methoxychlor	<0.50	<0.50	<0.50	<0.50	<.5	<0.50
Toxaphene	<5.0	<5.0	<5.0	<5.0	<.5	<5.0
Dimethoate	<.95	<5	<5	<.25	<5	<5.0
Disulfoton	<.95	<5	<5.0		<5	<5.0
Methyl parathion	<.95	<5	<5.0	<.25	<5	<5
Parathion	<.95	<5	<5.0	<.25	<5	<5.0
Phorate						
Thionazin	U	<5		<.25	<5	<5.0
2,4,5-T	0.055J	0.19J	<0.25	<0.25	0.12 J	<0.25
2,4,5-TP (Silvex)	<0.25	<0.25	<0.25	0.16J	<.25	0.12J
2,4-D	<0.50	1.4	1.7	1.0	1.3	1.4
Dinoseb	0.14J	0.16J	0.30	0.43	<.2	<0.20
1,1,1,2-Tetrachloroethane	<1.0	<1.0	<1.0	<1	<1	<1.0
1,1,1-Trichloroethane	<1.0	<1.0	<1.0	<1.0	<1	<1.0
1,1,2,2-Tetrachloroethane	<1.0	<1.0	<1.0	<1.0	<1	<1.0
1,1,2-Trichloroethane	<1.0	<1.0	<1.0	<1.0	<1	<1.0
1,1-Dichloroethane	<1.0	<1.0	<1.0	<1.0	<1	<1.0
1,1-Dichloroethene	<1.0	<1.0	<1.0	<1.0	<1	<1.0
1,1-Dichloropropene	<1.0	<1.0	<1.0	<1.0	<1	<1.0
1,2,3-Trichloropropane	<1.0	<1.0	<1.0	<1.0	<1	<1.0
1,2-Dibromo-3-chloropropane	<1.0	<1.0	<1.0	<1.0	<1	<1.0
1,2-Dibromoethane	<1.0	<1.0	<1.0	<1.0	<1	<1.0
1,2-Dichlorobenzene	<1.0	<1.0	<1.0	<1.0	<1	<1.0

Cell7 PLCRS

CELL 7 PLCRS						
	Aug_18	Dec_18	Jun_19	Dec_19	June_20	Dec_20
1,2-Dichloroethane	<1.0	<1.0	<1.0	<1.0	<1	<1.0
1,2-Dichloropropane	<1.0	<1.0	<1.0	<1.0	<1	<1.0
1,3-Dichlorobenzene	<1.0	<1.0	<1.0	<1.0	<1	<1.0
1,3-Dichloropropane	<1.0	<1.0	<1.0	<1.0	<1	<1.0
1,4-Dichlorobenzene	<1.0	<1.0	<1.0	<1.0	<1	<1.0
1,4-Dioxane (p-Dioxane)	0.59	2.7	<100 SIM 2.4ug/l	4.2	<100 SIM 3.3 ug/l	1.7
2,2-Dichloropropane	<1.0	<1.0	<1.0	<1.0	<1	<1.0
2-Butanone	16.7	14.4	10.8	13.1	14.2	3.6J
2-Hexanone	<5.0	<5.0	<5.0	<5.0	<5	<5.0
4-Methyl-2-pentanone	1.8J	1.6J	1.4J	<5.0	<5	<5.0
Acetone	274 D	195	103	179	124	49.7
Acetonitrile	62.9	156	128	193	<5	<5.0
Acrolein	<1.0	<1.0	<1.0	<1.0	<1	<1.0
Acrylonitrile	<1.0	<1.0	<1.0	<1.0	<1	<1.0
Allyl Chloride	<1.0	<1.0	<1.0	<1.0	<4	<4.0
Benzene	<1.0	<1.0	<1.0	<1.0	<1	<1.0
Bromochloromethane	<1.0	<1.0	<1.0	<1.0	<1	<1.0
Bromodichloromethane	<1.0	<1.0	<1.0	<1.0	<1	<1.0
Bromoform	<1.0	<1.0	<1.0	<1.0	<1	<1.0
Bromomethane	<1.0	<1.0	<1.0	<1.0	<1	<1.0
Carbon disulfide	<1.0	1.1	<1.0	<1.0	<1	1.1
Carbon tetrachloride	<1.0	<1.0	<1.0	<1.0	<1	<1.0
Chlorobenzene	<1.0	<1.0	<1.0	<1.0	<1	<1.0
Chloroethane	<1.0	<1.0	<1.0	<1.0	<1	<1.0
Chloroform	<1.0	<1.0	<1.0	<1.0	<1	<1.0
Chloromethane	<1.0	<1.0	<1.0	<1.0	<1	<1.0
Chloroprene	<1.0	<1.0	<1.0	<1.0	<1	<1.0
cis-1,2-Dichloroethene	<1.0	<1.0	<1.0	<1.0	<1	<1.0
cis-1,3-Dichloropropene	<1.0	<1.0	<1.0	<1.0	<1	<1.0
Dibromochloromethane	<1.0	<1.0	<1.0	<1.0	<1	<1.0
Dibromomethane	<1.0	<1.0	<1.0	<1.0	<1	<1.0
Dichlorodifluoromethane	<1.0	<1.0	<1.0	<1.0	<1	<1.0
Ethyl Methacrylate	<1.0	<1.0	<1.0	<1.0	<1	<1.0
Ethylbenzene	<1.0	<1.0	<1.0	<1.0	<1	<1.0
Iodomethane	<1.0	<1.0	<1.0	<1.0	4.2	<4.0
Isobutyl alcohol				5.8JJ	<20	<20.0
Methacrylonitrile	<1.0	<1.0	<1.0	<1.0	<1	<1.0
Methyl Methacrylate	<1.0	<1.0	<1.0	<1.0	<1	<1.0
Methylene chloride	<1.0	<1.0	<1.0	<1.0	<1	<1.0
Propionitrile	<4.0	<4.0	<4.0	<4.0	<4	<4.0
Silane, methoxytrimethyl-			<1.0			
Silanol, trimethyl-						
Styrene	<1.0	<1.0	<1.0		<1	<1.0
Tetrachloroethene	<1.0	<1.0	<1.0	<1.0	<1	<1.0
Toluene	<1.0	<1.0	<1.0	<1.0	<1	<1.0
trans-1,2-Dichloroethene	<1.0	<1.0	<1.0	<1.0	<1	<1.0
trans-1,3-Dichloropropene	<1.0	<1.0	<1.0	<1.0	<1	<1.0
trans-1,4-Dichloro-2-butene	<1.0	<1.0	<1.0	<1.0	<1	<1.0
Trichloroethene	<1.0	<1.0	<1.0	<1.0	<1	<1.0

Cell7 PLCRS

CELL 7 PLCRS						
	Aug_18	Dec_18	Jun_19	Dec_19	June_20	Dec_20
Trichlorofluoromethane	<1.0	<1.0	<1.0	<1.0	<1	<1.0
Trimethylsilyl fluoride+Sulfur diox						
Vinyl acetate	<1.0	<1.0	<1.0	<1.0	<1	<1.0
Vinyl chloride	<1.0	<1.0	<1.0	<1.0	<1	<1.0
Xylene (total)	<3.0	<3.0	<3.0	<3.0	<3	<3.0
1,2,4,5-Tetrachlorobenzene	U	<5.0	<5.0	<25.0 D	<5	<5.0
1,2,4-Trichlorobenzene	U	<5.0	<5.0	<25.0 D	<5	<5.0
1,2-Dichlorobenzene	U	<5.0	<5.0	<25.0 D	<5	<5.0
1,3,5-Trinitrobenzene	U	<5.0	<5.0	<25.0 D	<5	<5.0
1,3-Dichlorobenzene	U	<5.0	<5.0	<25.0 D	<5	<5.0
1,3-Dinitrobenzene	U	<5.0	<5.0	<25.0 D	<5	<5.0
1,4-Dichlorobenzene	U	<5.0	<5.0	<25.0 D	<5	<5.0
1,4-Naphthoquinone	U	<5.0	<5.0	<25.0 D	<5	<5.0
1-Naphthylamine	U	<5.0	<5.0	<25.0 D	<5	<5.0
2,2'-oxybis(1-chloropropane)		<5.0	<5.0	<25.0 D	<5	<5.0
2,3,4,6-Tetrachlorophenol	U	<5.0	<5.0	<25.0 D	<5	<5.0
2,4,5-Trichlorophenol	U	<5.0	<5.0	<25.0 D	<5	<5.0
2,4,6-Trichlorophenol	U	<5.0	<5.0	<25.0 D	<5	<5.0
2,4-Dichlorophenol	U	<5.0	<5.0	<25.0 D	<5	<5.0
2,4-Dimethylphenol	U	<5.0	<5.0	<25.0 D	<5	<5.0
2,4-Dinitrophenol	U	<10.0	<10.0	<50.0 D	<10	<10.0
2,4-Dinitrotoluene	U	<5.0	<5.0	<25.0 D	<5	<5.0
2,6-Dichlorophenol	U	<5.0	<5.0	<25.0 D	<5	<5.0
2,6-Dinitrotoluene	U	<5.0	<5.0	<25.0 D	<5	<5.0
2-Acetylaminofluorene	U	<5.0	<5.0	<25.0 D	<5	<5.0
2-Chloronaphthalene	U	<5.0	<5.0	<25.0 D	<5	<5.0
2-Chlorophenol	U	<5.0	<5.0	<25.0 D	<5	<5.0
2-Methylnaphthalene	U	<5.0	<5.0	<25.0 D	<5	<5.0
2-Methylphenol	0.328	<5.0	1.0J	<25.0 D	0.63 J	<5.0
2-Naphthylamine	U	<5.0	<5.0	<25.0 D	<5	<5.0
2-Nitroaniline	U	<5.0	<5.0	<25.0 D	<5	<5.0
2-Nitrophenol	U	<5.0	<5.0	<25.0 D	<5	<5.0
3,3'-Dichlorobenzidine	U	<5.0	<5.0	<25.0 D	<5	<5.0
3,3'-Dimethylbenzidine	U	<5.0	<5.0	<25.0 D	<5	<5.0
3-Methylcholanthrene	U	<5.0	<5.0	<25.0 D	<5	<5.0
3-Methylphenol/4-Methylphenol	46.8	39.1	110 D		44.4	1.2J
3-Nitroaniline	U	<5.0	<5.0	<25.0 D	<5	<5.0
4,6-Dinitro-2-methylphenol	U	<10.0	<10.0	<50.0 D	<10	<10.0
4-Aminobiphenyl	U	<5.0	<5.0	<25.0 D	<5	<5.0
4-Bromophenyl-phenylether	U	<5.0	<5.0	<25.0 D	<5	<5.0
4-Chloro-3-methylphenol	U	<5.0	<5.0	<25.0 D	<5	<5.0
4-Chloroaniline	U	<5.0	<5.0	<25.0 D	<5	<5.0
4-Chlorophenyl-phenylether	U	<5.0	<5.0	<25.0 D	<5	<5.0
4-Nitroaniline	U	<5.0	<5.0	<25.0 D	<5	<5.0
4-Nitrophenol	U	<10.0	<10.0	<50.0 D	<10	<10.0
5-Nitro-o-toluidine	U	<5.0	<5.0	<25.0 D	<5	<5.0
7,12-Dimethylbenz(a)anthracene		<5.0	<5.0	<25.0 D	<5	<5.0
Acenaphthene	U	<5.0	<5.0	<25.0 D	<5	<5.0
Acenaphthylene	U	<5.0	<5.0	<25.0 D	<5	<5.0

Cell7 PLCRS

CELL 7 PLCRS						
	Aug_18	Dec_18	Jun_19	Dec_19	June_20	Dec_20
Acetophenone	U	<5.0	<5.0	<25.0 D	<5	<5.0
Anthracene	U	<5.0	<5.0	<25.0 D	<5	<5.0
Benzo(a)anthracene	U	<5.0	<5.0	<25.0 D	<5	<5.0
Benzo(a)pyrene	U	<5.0	<5.0	<25.0 D	<5	<5.0
Benzo(b)fluoranthene	U	<5.0	<5.0	<25.0 D	<5	<5.0
Benzo(g,h,i)perylene	U	<5.0	<5.0	<25.0 D	<5	<5.0
Benzo(k)fluoranthene	U	<5.0	<5.0	<25.0 D	<5	<5.0
Benzyl alcohol	U	<5.0	<5.0	<25.0 D	<5	0.88J
Bis(2-chloroethoxy)methane	U	<5.0	<5.0	<25.0 D	<5	<5.0
Bis(2-chloroethyl)ether	U	<5.0	<5.0	<25.0 D	<5	<5.0
Bis(2-ethylhexyl)phthalate	U	<5.0	<5.0	8.9J D	<5	<5.0
Butyl benzyl phthalate	U	<5.0	<5.0	<25.0 D	<5	<5.0
Chlorobenzilate	U	<5.0	<5.0	<25.0 D	<5	<5.0
Chrysene	U	<5.0	<5.0	<25.0 D	<5	<5.0
Diallate	U	<5.0	<5.0	<25.0 D	<5	<5.0
Dibenzo(a,h)anthracene	U	<5.0	<5.0	<25.0 D	<5	<5.0
Dibenzofuran	U	<5.0	<5.0	<25.0 D	<5	<5.0
Diethylphthalate	U	<5.0	<5.0	<25.0 D	<5	<5.0
Dimethylphthalate	U	<5.0	<5.0	<25.0 D	<5	<5.0
Di-n-butyl phthalate	U	<5.0	<5.0	<25.0 D	<5	<5.0
Di-n-octyl phthalate	U	<5.0	<5.0	<25.0 D	<5	<5.0
Ethyl methanesulfonate	U	<5.0	<5.0	<25.0 D	<5	<5.0
Famphur	<.95	<10.0	<10.0	<50.0 D	<10	<10.0
Fluoranthene	U	<5.0	<5.0	<25.0 D	<5	<5.0
Fluorene	U	<5.0	<5.0	<25.0 D	<5	<5.0
Hexachlorobenzene	U	<5.0	<5.0	<.25	<5	<5.0
Hexachlorobutadiene	U	<5	<5	<025	<5	<5.0
Hexachlorocyclopentadiene	U	<5	<5.0	<25.0 D	<5	<5.0
Hexachloroethane	U	<5.0	<5.0	<25.0 D	<5	<5.0
Hexachloropropene	U	<5.0	<5.0	<25.0 D	<5	<5.0
Indeno(1,2,3-cd)pyrene	U	<5.0	<5.0	<25.0 D	<5	<5.0
Isodrin	U	<5.0	<5.0	<25.0 D	<5	<5.0
Isophorone	U	<5.0	<5.0	<25.0 D	<5	<5.0
Isosafrole	U	<5.0	<5.0	<25.0 D	<5	<5.0
Kepone	U	<10.0	<10.0	<50.0 D	<10	<10.0
Methapyrilene	U	<5.0	<5.0	<25.0 D	<5	<5.0
Methyl methanesulfonate	U	<5		<25.0 D	<5	<5.0
Naphthalene	U	<5.0	<5.0	<25.0 D	<5	<5.0
Nitrobenzene	U	<5.0	<5.0	<25.0 D	<5	<5.0
N-Nitrosodiethylamine	U	<5	<5.0	<25.0 D	<5	<5.0
N-Nitrosodimethylamine	U	<5	<5.0	<25.0 D	<5	<5.0
N-Nitroso-di-n-butylamine	U	<5.0	<5.0	<25.0 D	<5	<5.0
N-Nitroso-di-n-propylamine	U	<5.0	<5.0	<25.0 D	<5	<5.0
N-Nitrosodiphenylamine	U	<5.0	<5.0	<25.0 D	<5	<5.0
N-Nitrosomethylethylamine	U	<5.0	<5.0	<25.0 D	<5	<5.0
N-Nitrosopiperidine	U	<5.0	<5.0	<25.0 D	<5	<5.0
N-Nitrosopyrrolidine	U	<5.0	<5.0	<25.0 D	<5	<5.0
O,O,O-Triethylphosphorothioate	U	<5.0	<5.0	<25.0 D	<5	<5.0
o-Toluidine	U	<5.0	<5.0	<25.0 D	<5	<5.0

Cell7 PLCRS

CELL 7 PLCRS						
	Aug_18	Dec_18	Jun_19	Dec_19	June_20	Dec_20
p-Dimethylaminoazobenzene	U	<5.0	<5.0	<25.0 D	<5	<5.0
Pentachlorobenzene	U	<5.0	<5.0	<25.0 D	<5	<5.0
Pentachloronitrobenzene	U	<5.0	<5.0	<25.0 D	<5	<5.0
Pentachlorophenol	2.37	<10.0	<10.0	<50.0 D	<10	<10.0
Phenacetin	U	<5.0	<5.0	<25.0 D	<5	<5.0
Phenanthrene	U	<5.0	<5.0	<25.0 D	<5	<5.0
Phenol	52.2	31.4	115 D	70.0 D	87.1	3.1J
p-Phenylenediamine	U	<5	<10.0	<50	<10	<10.0
Pronamide	U	<5.0	<5.0	<.25	<5	<5.0
Sulfide	1.6J	8	8.0	4.8	25.6	<2.0
2378-TCDF	ND	ND		ND	ND	ND
2378-TCDD	ND	ND	ND	ND	ND	ND
Total Uranium	0.347 ± 0.013 (0.262) C:NA T:NA	.855±.049 (2.62) C:NA T:NA	0.281 ± 0.014 (0.262) C:NA T:NA	0.789 ± 0.039 (0.262) C:NA T:NA	0.751 ± 0.045 (2.620) C:NA T:NA	0.526 ± 0.049 (2.620) C:NA T:NA
Perfluorobutanesulfonic acid PFBS	130	130		170	160	120
Perfluoroheptanoic acid PFHpA	19	18		24	26	35
Perfluorohexanesulfonic acid PFHxS	4.7	4.2		11B	8.6	5.9
Perfluorononanoic acid PFNA	1.7	1.2		1.4J	5	2
Perfluorooctanesulfonic acid PFOS	3.3	2		3	16	4
Perfluorooctanoic acid PFOA	22	22		32	50	47
Radium-226	6.34 ± 2.29 (1.80) C:NA T:42%	15.7 ± 7.46 (2.36) C:NA T:88%	9.05 ± 2.77 (0.511) C:NA T:85%	2.93 ± 1.62 (1.44) C:NA T:61%	3.77 ± 2.18 (0.852) C:NA T:43%	1.21 ± 0.852 (0.938) C:NA T:89%
Radium-228	10.2 ± 3.75 (5.39) C:72% T:85%	6.62 ± 2.38 (3.68) C:80% T:89%	6.45 ± 1.59 (1.46) C:78% T:52%	3.90 ± 2.48 (4.69) C:81% T:24%	7.79 ± 2.29 (2.88) C:78% T:33%	3.50 ± 1.31 (2.03) C:79% T:37%
6:2 FTS	5.4	6.6		11J	10	6.2
8:2 FTS	19U	ND		ND	ND	ND
N-ethyl perfluorooctansulfamidoacetic acidNEtFOSAA	19U	ND		ND	ND	ND
N-methylperfluorooctansulfamicacetic acid NMeFOSAA	19U	ND		ND	ND	ND
perfluorobutanoic acid PFBA	260	170		180	260	170
perfluorodecansulfonic acid PFDS	19U	ND		ND	ND	ND
perfluorodecanoic acid PFDA	4.5	0.44		.38J	3.2	0.55
perfluorododecanoic acid PFDoA	19U	ND		ND	ND	ND
perfluoroheptanesulfonic acid PFHps	19U	ND		ND	0.26	ND
perfluorohexanoic acid PFHxA	210	250		320	370	350
perfluorooctane sulfonamide FOSA	19U	ND		1J	2.2	ND
perfluoropentanoic acid PFPeA	100	94		130	140	120
perfluorotetradecanoic acid PFTeA	19U	ND		ND	ND	ND
perfluorotridecnaoic acid PFTriA	19U	ND		ND	ND	ND
perfluoroundecanoic acid PFUnA	19U	ND		ND	ND	ND
n-Nitrosomorpholine	U					
Dimethylbenz(A) Anthracene	U					
Bis(2-chloroisopropyl)ether	U					
total PFOA/PFAS	760.6	698.44		859	1051.26	860.65



## Appendix 1

December 2020 Pace Analytical Laboratory Report and QA/QC

(see attached CD)

# BABYLON LANDFILL - FIELD DATA - DECEMBER 10, 2020

## Traditional Wells - Groundwater Sampling Data

WELL #	Well Survey Elevation	Well Size	Metal or PVC	TPVC (in ft) (Top of PVC)	TOC (in ft) (Top of Casing)	BOC (in ft) (Bottom of Casing)	One Well Volume (Gallons)	Three Well Volumes (Gallons)	Groundwater Contour Levels
GM-2D	69.25	4"	PVC	26.54	27.35	86.00	38.30	114.90	41.90
GM-4D	62.43	4"	PVC	19.30	19.95	91.40	46.66	139.97	42.48
GM-5D	62.35	4"	PVC	18.73	20.15	91.80	46.79	140.36	42.20
GM-6D	63.84	4"	PVC	21.23	21.40	92.80	46.62	139.87	42.44
GM-7D	63.23	4"	PVC	20.25	20.93	91.10	45.82	137.46	42.30
GM-15D	50.74	4"	PVC	12.58	13.04	84.50	46.66	139.99	37.70
GM-16D	?	4"	PVC	9.07	9.42	87.00	50.66	151.98	?
GM-17D	52.09	4"	PVC	14.42	14.81	87.70	47.60	142.79	37.28
GM-18D	?	4"	PVC	14.92	15.32	78.00	40.93	122.79	?
GM-19D	53.34	4"	PVC	14.73	14.94	87.40	47.32	141.95	38.40

WELL #	Start Purge	Stop Purge	Well Notes For Sampling
GM-2D	820	900	Clear, no odors
GM-4D	1216	1248	Clear, no odors
GM-5D	1218	1257	Clear, yellow tint, no odors
GM-6D	1227	1305	Cloudy, no odors
GM-7D	1300	1322	Clear, small black particles, no odors
GM-15D	1040	1110	Clear, foamy, no odors
GM-16D	1045	1118	Clear, small orange particles, no odors
GM-17D	1130	1205	Clear, no odors
GM-18D	956	1030	Clear, some particles, no odors
GM-19D	915	948	Clear, no odors

Water Quality Parameters									
WELL #	Sampling Date	Sample Time	pH (SU)	ORP (mv)	Conductivity (umhos/cm2)	Temp. (oC)	Turbidity (NTU)	Dis. Oxygen (DO) mg/L	
GM-2D	12/10/2020	905	7.12	-62.2	187.9	14.1	11.00	1.92	
GM-4D	12/10/2020	1250	7.31	-67.6	259	14.8	4.60	4.34	
GM-5D	12/10/2020	1300	7.95	-102.1	566	15.9	50.00	3.35	
GM-6D	12/10/2020	1310	8.46	-128.7	844	16.3	31.00	4.51	
GM-7D	12/10/2020	1326	9.16	-166.3	682	17.1	18.00	6.79	
GM-15D	12/10/2020	1112	8.68	-140.0	1089	14.2	12.00	3.37	
GM-16D	12/10/2020	1120	8.39	-123.1	300	16.7	16.00	2.54	
GM-17D	12/10/2020	1207	7.41	-75.5	202	16.9	7.70	3.04	
GM-18D	12/10/2020	1032	8.39	-127.8	628	12.5	8.50	3.05	
GM-19D	12/10/2020	948	7.70	-58.9	234	13.0	4.60	5.97	

# BABYLON LANDFILL - FIELD DATA - DECEMBER 10, 2020

## Leachate Sampling Data

WELL #	Date	Start Purge	Stop Purge	Gallons Purged	Well Notes For Sampling
NNU-PLCRS	12/10/2020	1415	1418	~ 40	Yellow tint, black particles, odors, sample warm
NNU-SLCRS	12/10/2020	1407	1409	~ 40	Clear, odors, sample warm
ONU-SLCRS	12/10/2020	1350	1352	~ 60	Clear, odors
SA-SLCRS	12/10/2020	Direct Sample	Direct Sample	0	Cloudy to turbid, black in color, no odors
CELL - 7	12/10/2020	Direct Sample	Direct Sample	0	Clear, odors

## Leachate Parameters

WELL #	Sampling Time	pH (SU)	ORP (mv)	Conductivity (umhos/cm2)	Temp. (oC)	Turbidity (NTU)	Dissolved Oxygen (DO) mg/L
NNU-PLCRS	1420	8.56	-139.0	>20,000	21.7	6.90	0.34
NNU-SLCRS	1410	7.37	-72.0	>20,000	28.6	2.30	0.26
ONU-SLCRS	1355	8.02	-106.6	>20,000	14.1	15.00	4.92
SA-SLCRS	1435	8.59	-132.8	1,200	15.6	70.00	4.05
CELL - 7	1454	7.93	-102.4	>20,000	15.8	5.50	2.02

**Field Notes:**

NNU-PLCRS: **New Northern U Primary** \* One Tap Location for Primary/Secondary (Top Road)

NNU-SLCRS: **New Northern U Secondary** \* One Tap Location for Primary/Secondary (Top Road)

ONU-SLCRS: **Old Northern U Secondary** \*One Tap Location for Primary/Secondary (Lower Road)

SA-SLCRS: **Southern Ash Secondary** \*Use Bailer / Square Metal Door

CELL 7: **Primary System** \* Use Bailer / First Round Black Cover (Left Cover)

# BABYLON LANDFILL - FIELD DATA - DECEMBER 11, 2020

## Wells GM-26 to GM-28 / Groundwater Sampling Data

WELL #	Well Survey Elevation	Well Size	Metal or PVC	TPVC (in ft) (Top of PVC)	TOC (in ft) (Top of Casing)	BOC (in ft) (Bottom of Casing)	One Well Volume (Gallons)	Three Well Volumes (Gallons)	Groundwater Contour Levels
GM-26		4"	*PVC	20.36	*	32.50	7.93	23.78	
GM-26I		4"	*PVC	20.00	*	42.50	14.69	44.08	
GM-27		4"	PVC	26.21	26.57	36.70	6.61	19.84	
GM-27I		4"	PVC	26.51	26.74	47.50	13.56	40.67	
GM-28		4"	PVC	25.93	26.12	37.50	7.43	22.29	
GM-28I		4"	PVC	26.15	26.32	46.91	13.45	40.34	

### Well Notes For Sampling

WELL #	Start Purge	Stop Purge	Well Notes For Sampling
GM-26	815	902	Cloudy, orange tint, no odors
GM-26I	813	856	Cloudy, orange tint, no odors
GM-27	918	938	Clear, yellow tint, no odors
GM-27I	920	950	Turbid, grey to black in color, odor
GM-28	1018	1048	Cloudy to slightly turbid, no odors
GM-28I	1020	1100	Clear, a lot of small black particles, slight odors

### Water Quality Parameters

WELL #	Sampling Date	Sample Time	pH (SU)	ORP (mv)	Conductivity (umhos/cm2)	Temp. (oC)	Turbidity (NTU)	Dis. Oxygen (DO) mg/L
GM-26	12/11/2020	908	7.65	-87.9	464	17.7	850.0	4.41
GM-26I	12/11/2020	900	7.11	-59.8	340	14.9	290.0	4.06
GM-27	12/11/2020	1000	8.96	-154.3	1082	14.3	20.0	3.38
GM-27I	12/11/2020	952	8.80	-146.9	1106	13.1	750.0	2.71
GM-28	12/11/2020	1052	8.16	-114.7	2060	15.2	95.0	3.04
GM-28I	12/11/2020	1103	9.62	-190.5	645	14.5	18.0	2.23

Field Notes: Duplicate (GM-X) performed on GM-27I @ 954  
 Equipment Blank @ 935 w/new bailer  
 MS/MSD performed on GM-28I @ 1105  
**GM-28I did not have a cover to the flush mount well**

- Notes: N/F : Not found due to high grass or deep snow.  
 N/S : No sample due to dry well or frozen well from extreme cold temps.  
 \*PVC ABOVE TOC

## PFCs Sampling Checklist

Date: Thu. 12-10-2020

Weather (temp./precipitation): Sunny 50° Site Name: Babylon Landfill

### **Field Clothing and PPE:**

- No clothing or boots containing Gore-Tex™
- All safety boots made from polyurethane and PVC
- No materials containing Tyvek®
- Field crew has not used fabric softener on clothing
- Field crew has not used cosmetics, moisturizers, hand cream, or other related products this morning
- Field crew has not applied unauthorized sunscreen or insect repellent

### **Field Equipment:**

- No Teflon® or LDPE containing materials on-site
- All sample materials made from stainless steel, HDPE, acetate, silicon, or polypropylene
- No waterproof field books on-site
- No plastic clipboards, binders, or spiral hard cover notebooks on-site
- No adhesives (Post-It Notes) on-site

- Coolers filled with regular ice only. No chemical (blue) ice packs in possession

### **Sample Containers:**

- All sample containers made of HDPE or polypropylene
- Caps are unlined and made of HDPE or polypropylene

### **Wet Weather (as applicable):**

- Wet weather gear made of polyurethane and PVC only

### **Equipment Decontamination:**

- "PFC-free" water on-site for decontamination of sample equipment. No other water sources to be used.
- Alconox and Liquinox to be used as decontamination materials

### **Food Considerations:**

- No food or drink on-site with exception of bottled water and/or hydration drinks (i.e., Gatorade and Powerade) that is available for consumption only in the staging area

If any applicable boxes cannot be checked, the Field Lead shall describe the noncompliance issues below and work with field personnel to address noncompliance issues prior to commencement of that day's work. Corrective action shall include removal of noncompliance items from the site or removal of worker offsite until in compliance.

Describe the noncompliance issues (include personnel not in compliance) and action/outcome of noncompliance:

Field Lead Name: Brian Nichols

Field Lead Signature: Brian Time: 800 AM

# PFCs Sampling Checklist

Date: Fri. 12-11-2020

Weather (temp./precipitation): Sunny 53° Site Name: Babylon Landfill

## Field Clothing and PPE:

- No clothing or boots containing Gore-Tex™
- All safety boots made from polyurethane and PVC
- No materials containing Tyvek®
- Field crew has not used fabric softener on clothing
- Field crew has not used cosmetics, moisturizers, hand cream, or other related products this morning
- Field crew has not applied unauthorized sunscreen or insect repellent

## Field Equipment:

- No Teflon® or LDPE containing materials on-site
- All sample materials made from stainless steel, HDPE, acetate, silicon, or polypropylene
- No waterproof field books on-site
- No plastic clipboards, binders, or spiral hard cover notebooks on-site
- No adhesives (Post-It Notes) on-site

- Coolers filled with regular ice only. No chemical (blue) ice packs in possession

## Sample Containers:

- All sample containers made of HDPE or polypropylene
- Caps are unlined and made of HDPE or polypropylene

## Wet Weather (as applicable):

- Wet weather gear made of polyurethane and PVC only

## Equipment Decontamination:

- "PFC-free" water on-site for decontamination of sample equipment. No other water sources to be used.
- Alconox and Liquinox to be used as decontamination materials

## Food Considerations:

- No food or drink on-site with exception of bottled water and/or hydration drinks (i.e., Gatorade and Powerade) that is available for consumption only in the staging area

If any applicable boxes cannot be checked, the Field Lead shall describe the noncompliance issues below and work with field personnel to address noncompliance issues prior to commencement of that day's work. Corrective action shall include removal of noncompliance items from the site or removal of worker offsite until in compliance.

Describe the noncompliance issues (include personnel not in compliance) and action/outcome of noncompliance:

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Field Lead Name: Brian Nichols

Field Lead Signature: Brian Nichols Time: 800 AM

## Appendix 2

### Baseline and Expanded Parameters List (6NYCRR Part 363-4.6(h))

(5) Data quality assessment. At the conclusion of each sampling event and analysis of the samples collected, data quality assessment must occur. A data quality assessment report must be submitted with the results from each sampling event. Data quality assessment must occur in two phases – data validation and data usability analysis.

(i) Data validation.

(a) For those sampling events for which only routine parameters are analyzed, the required data validation may be performed by the laboratory that performed the sample analyses.

(b) For those sampling events in which groundwater samples are analyzed for baseline or expanded parameters, the data validation must be performed by a person with experience with similar validation projects and who is not affiliated with the laboratory that performed the analyses and who is acceptable to the department.

(c) The data validation must be performed on all analytical data for the facility at a rate acceptable to the department, but not less than five percent of the data generated, and must consist, at a minimum, of the following:

(1) field records and analytical data are reviewed to determine whether the data are accurate and defensible. All AQA/AQC information must be reviewed along with any corrective actions taken during that sampling event, and

(2) all data summaries must be clearly marked to identify any data that are not representative of environmental conditions at the site, or that were not generated in accordance with the site analytical plan.

(ii) Data usability analysis.

(a) The data usability analysis must be performed on all analytical data generated by the requirements for this Part for the facility and must consist of the following:

(1) an assessment to determine if the data quality objectives were met;

(2) for consistency, comparison of the analytical data with the results from previous sampling events;

(3) evaluation of field duplicate results to indicate the samples are representative;

(4) comparison of the results of all field blanks, trip blanks, equipment rinse blanks, and method blanks with full data sets to provide information concerning contaminants that may have been introduced during sampling, shipping, or analysis;

(5) evaluation of matrix effects to assess the performance of the analytical method with respect to the sample matrix, and determine whether the data have been biased high or low due to matrix effects;

(6) integration of the field and laboratory data with geological, hydrogeological, and meteorological data to provide information about the extent of contamination, if it occurs; and

(7) comparison of precision, accuracy, representativeness, comparability, completeness, and defensibility of the data generated with that required to meet the data quality objectives established in the site analytical plan.

(h) Water quality analysis tables.

The water quality analysis tables in this section list the routine, baseline, and expanded parameters for analysis of all monitoring samples. The department may modify the parameters for analysis based on the location of the landfill or site-specific characteristics of waste disposed at the landfill.

TABLE 1: ROUTINE PARAMETERS <sup>1</sup>

Common Name (and CAS number, as appropriate) <sup>2</sup>		
Field Parameters	Leachate Indicators:	Inorganic Parameters (total)
Static water level (in wells and sumps)	Total Kjeldahl Nitrogen	Arsenic
Specific Conductance	Ammonia (7664-41-7)	Cadmium
Temperature	Nitrate	Calcium
Floaters or Sinkers <sup>3</sup>	Chemical Oxygen Demand	Iron
Temperature	Biochemical Oxygen Demand (BOD <sub>5</sub> )	Lead
pH	Total Organic Carbon	Magnesium
Eh	Total Dissolved Solids	Manganese
Dissolved Oxygen <sup>4</sup>	Sulfate	Potassium
Field Observations <sup>5</sup>	Alkalinity	Sodium
Turbidity	Phenols (108-95-2)	
	Chloride	
	Bromide (24959-67-9)	
	Total hardness as CaCO <sub>3</sub>	

TABLE 2A: BASELINE PARAMETERS: Field Parameters, Leachate Indicators, and Inorganic Parameters <sup>6</sup>



Common Name (and CAS number, as appropriate) <sup>7</sup>		
Field Parameters:	Leachate Indicators:	Inorganic Parameters (total unless otherwise noted):
Static water level (in wells and sumps)	Total Kjeldahl Nitrogen	Aluminum
Specific Conductance	Ammonia (7664-41-7)	Antimony
Temperature	Nitrate	Arsenic
Floaters or Sinkers <sup>8</sup>	Chemical Oxygen Demand	Barium
Temperature	Biochemical Oxygen Demand (BOD <sub>5</sub> )	Beryllium
pH	Total Organic Carbon	Cadmium
Eh	Total Dissolved Solids	Calcium
Dissolved Oxygen <sup>9</sup>	Sulfate	Chromium
Field Observations <sup>10</sup>	Alkalinity	Chromium (Hexavalent) <sup>11</sup>
Turbidity	Phenols (108-95-2)	Cobalt
	Chloride	Copper
	Bromide (24959-67-9)	Cyanide
	Total hardness as CaCO <sub>3</sub>	Iron
	Color	Lead
	Boron (7440-42-8)	Magnesium
		Manganese
		Mercury
		Nickel
		Potassium
		Selenium
		Silver
		Sodium
		Thallium
		Vanadium
		Zinc

TABLE 2B: BASELINE PARAMETERS: Organic Parameters<sup>12</sup>

Common Name (and CAS number, as appropriate) <sup>13</sup>		
Organic Parameters:		
Acetone (67-64-1)	1,1-Dichloroethane; Ethylidene chloride (75-34-3)	Styrene (100-42-5)
Acrylonitrile (107-13-1)	1,2-Dichloroethane; Ethylene dichloride (107-06-02)	1,1,1,2-Tetrachloroethane (630-20-6)
Benzene (71-43-2)	1,1-Dichloroethylene; 1,1-Dichloroethene; Vinylidene chloride (75-35-4)	1,1,2,2-Tetrachloroethane (79-34-5)
Bromochloromethane (74-97-5)	cis-1,2-Dichloroethylene; cis-1,2-Dichloroethene (156-59-2)	Tetrachloroethylene; Tetrachloroethene; Perchloroethylene (127-18-4)
Bromodichloromethane (75-27-4)	trans-1,2-Dichloroethylene; trans-1,2-Dichloroethene (156-60-2)	Toluene (108-88-3)
Bromoform; Tribromomethane (75-25-2)	1,2-Dichloropropane; Propylene dichloride (78-87-5)	1,1,1-Trichloroethane; Methylchloroform (71-55-6)
Carbon disulfide (75-15-0)	cis-1,3-Dichloropropene (10061-01-5)	1,1,2-Trichloroethane (79-00-5)
Carbon tetrachloride (56-23-5)	trans-1,3-Dichloropropene (10061-02-6)	Trichloroethylene; Trichloroethene (79-01-6)
Chlorobenzene (108-90-7)	Ethylbenzene (100-41-4)	Trichlorofluoromethane; CFC-11 (75-69-4)
Chloroethane; Ethyl chloride (75-00-3)	2-Hexanone; Methyl butyl ketone (591-78-6)	1,2,3-Trichloropropane (96-18-4)
Chloroform; Trichloromethane (67-66-3)	Methyl bromide; Bromomethane (74-83-9)	Vinyl acetate (108-05-4)
Dibromochloromethane; Chlorodibromomethane (124-48-1)	Methyl chloride; Chloromethane (74-87-3)	Vinyl chloride; Chloroethene (75-01-4)
1,2-Dibromo-3-chloropropane; DBCP (96-12-8)	Methylene bromide; Dibromomethane (74-95-3)	Xylenes (1330-20-7)
1,2-Dibromoethane; Ethylene dibromide; EDB (106-93-4)	Methylene chloride; Dichloromethane (75-09-2)	
o-Dichlorobenzene; 1,2-Dichlorobenzene (95-50-1)	Methyl ethyl ketone; MEK; 2-Butanone (78-93-3)	
p-Dichlorobenzene; 1,4-Dichlorobenzene (106-46-7)	Methyl iodide; Iodomethane (74-88-4)	
trans-1,4-Dichloro-2-butene (110-57-6)	4-Methyl-2-pentanone; Methyl isobutyl ketone (108-10-1)	

TABLE 3A: EXPANDED PARAMETERS: Field Parameters, Leachate Indicators, Radionuclides, and Inorganic Parameters<sup>14</sup>

Common Name (and CAS number, as appropriate) <sup>15</sup>

Field Parameters:	Leachate Indicators:	Inorganic Parameters: (total unless otherwise noted)	Radionuclides <sup>16</sup>
Static water level (in wells and sumps)	Total Kjeldahl Nitrogen	Aluminum	Radium-226 per EPA 903.1
Specific Conductance	Ammonia (7664-41-7)	Antimony	Radium-228 per EPA 904.0
Temperature	Nitrate	Arsenic	Total Uranium per EPA 908.0
Floaters or Sinkers <sup>17</sup>	Chemical Oxygen Demand	Barium	
Temperature	Biochemical Oxygen Demand (BOD <sub>5</sub> )	Beryllium	
pH	Total Organic Carbon	Cadmium	
Eh	Total Dissolved Solids	Calcium	
Dissolved Oxygen <sup>18</sup>	Sulfate	Chromium	
Field Observations <sup>19</sup>	Alkalinity	Chromium (Hexavalent) <sup>20</sup>	
Turbidity	Phenols (108-95-2)	Cobalt	
	Chloride	Copper	
	Bromide (24959-67-9)	Cyanide	
	Total hardness as CaCO <sub>3</sub>	Iron	
	Color	Lead	
	Boron (7440-42-8)	Magnesium	
		Manganese	
		Mercury	
		Nickel	
		Potassium	
		Selenium	
		Silver	
		Sodium	
		Thallium	
		Tin	
		Vanadium	
		Zinc	

TABLE 3B: EXPANDED PARAMETERS: Organic Parameters<sup>21</sup>

Common Name (and CAS number, as appropriate) <sup>22</sup>		
Organic Parameters:		
Acenaphthene (83-32-9)	2,4-Dichlorophenol (120-83-2)	Naphthalene (91-20-3)
Acenaphthylene (208-96-8)	2,6-Dichlorophenol (87-65-0)	1,4-Naphthoquinone (130-15-4)
Acetone (67-64-1)	1,2-Dichloropropane; Propylene dichloride (78-87-5)	1-Naphthylamine (134-32-7)
Acetonitrile, Methyl cyanide (75-05-8)	1,3-Dichloropropane, Trimethylene dichloride (142-28-9)	2-Naphthylamine (91-59-8)
Acetophenone (98-86-2)	2,2-Dichloropropane, Isopropylidene chloride (594-20-7)	o-Nitroaniline, 2-Nitroaniline (88-74-4)
2-Acetylamino fluorene; 2-AAF (53-96-3)	1,1-Dichloropropene (563-58-6)	m-Nitroaniline; 3-Nitroaniline (99-09-2)
Acrolein (107-02-8)	cis-1,3-Dichloropropene (10061-01-5)	p-Nitroaniline, 4-Nitroaniline (100-01-6)
Acrylonitrile (107-13-1)	trans-1,3-Dichloropropene (10061-02-6)	Nitrobenzene (98-95-3)
Aldrin (309-00-2)	Dieldrin (60-57-1)	o-Nitrophenol 2-Nitrophenol (88-75-5)
Allyl chloride (107-05-1)	Diethyl phthalate (84-66-2)	p-Nitrophenol; 4-Nitrophenol (100-02-7)
4-aminobiphenyl (92-67-1)	0,0-Diethyl 0-2-pyrazinyl	N-Nitrosodi-n-butylamine (924-16-3)
Anthracene (120-12-7)	cis-1,2-Dichloroethylene; cis-1,2-Dichloroethene (156-59-2)	
N-Nitrosodiethylamine (55-18-5)		
Benzene (71-43-2)	trans-1,2-Dichloroethylene (156-60-2)	N-Nitrosodimethylamine (62-75-9)
Benzo[a]anthracene, Benzanthracene (56-55-3)	Phosphorothioate, Thionazin (297-97-2)	N-Nitrosodiphenylamine (86-30-6)
Benzo[b]fluoranthene (205-99-2)	Dimethoate (60-51-5)	N-Nitrosodipropylamine; N-Nitroso-N-dipropyl-amine, Di-n-propylnitrosamine (621-64-7)
Benzo[k]fluoranthene (207-08-9)	p-(Dimethylamino)azobenzene (60-11-7)	N-Nitrosomethylethylamine (10595-95-6)
Benzo[ghi]perylene (191-24-2)	7,12-Dimethylbenz[a]anthracene (57-97-6)	N-Nitrosopiperidine (100-75-4)
Benzo[a]pyrene (50-32-8)	3,3 <sup>21</sup> -Dimethylbenzidine (119-93-7)	N-Nitrosopyrrolidine (930-55-2)
Benzyl alcohol (100-51-6)	2,4-Dimethylphenol, m-Xylenol (105-67-9)	5-Nitro-o-toluidine (99-55-8)
alpha-BHC (319-84-6)	Dimethyl phthalate (131-11-3)	Parathion (56-38-2)
beta-BHC (319-85-7)	m-Dinitrobenzene (99-65-0)	Pentachlorobenzene (608-93-5)
delta-BHC (319-86-8)	4,6-Dinitro-o-cresol 4,6-Dinitro-2-methylphenol (534-52-1)	Pentachloronitrobenzene (82-68-8)

gamma-BHC, Lindane (58-89-9)	2,4-Dinitrophenol (51-28-5)	Pentachlorophenol (87-86-5)
Bis(2-chloroethoxy)methane (111-91-1)	2,4-Dinitrotoluene (121-14-2)	Phenacetin (62-44-2)
Bis(2-chloroethyl) ether, Dichloroethyl ether (111-44-4)	2,6-Dinitrotoluene (606-20-2)	Phenanthrene (85-01-8)
Bis-(2-chloro-1-methyl-ethyl)ether, 2,2 <sup>21</sup> -Dichlorodiisopropyl ether, DCIP <sup>23</sup>	Dinoseb, DNBP; 2-sec-Butyl-4,6-dinitrophenol (88-85-7)	Phenol (108-95-2)
Bis(2-ethylhexyl)phthalate (117-81-7)	Di-n-octyl phthalate (117-84-0)	p-Phenylenediamine (106-50-9)
Bromochloromethane (74-97-5)	Diphenylamine (122-39-4)	Phorate (298-02-2)
Bromodichloromethane (75-27-4)	Disulfoton (298-04-4)	Polychlorinated biphenyls; PCBs; Aroclors <sup>24</sup>
Bromoform (75-25-2)	Endosulfan I (959-98-8)	Polychlorinated dibenzo-p-dioxins; PCDDs <sup>25</sup>
4-Bromophenyl phenyl ether (101-55-3)	Endosulfan II (33213-65-9)	Polychlorinated dibenzo-furans; PCDFs <sup>26</sup>
Butyl benzyl phthalate, Benzyl butyl phthalate (117-81-7)	Endosulfan sulfate (1031-07-8)	Pronamide (23950-58-5)
Carbon disulfide (75-15-0)	Endrin (72-20-8)	Propionitrile; Ethyl cyanide (107-12-0)
Carbon tetrachloride (56-23-5)	Endrin aldehyde (7421-93-4)	Pyrene (129-00-0)
Chlordane <sup>27</sup>	Ethylbenzene (100-41-4)	Safrole (94-59-7)
p-Chloroaniline (106-47-8)	Ethyl methacrylate (97-63-2)	Silvex, 2,4,5-TP (93-72-1)
Chlorobenzene (108-90-7)	Ethyl methanesulfonate (62-50-0)	Styrene (100-42-5)
Chlorobenzilate (510-15-6)	Famphur (52-85-7)	2,4,5-T, 2,4,5-trichloro- phenoxyacetic acid (93-76-5)
p-Chloro-m-cresol; 4-Chloro-3-methylphenol (59-50-7)	Fluoranthene (206-44-0)	1,2,4,5-Tetrachlorobenzene (95-94-3)
Chloroethane, Ethyl chloride (75-00-3)	Fluorene (86-73-7)	2,3,7,8-Tetrachlorodi- benzo-p-dioxin, 2,3,7,8-TCDD (1746-01-6)
Chloroform; Trichloromethane (67-66-3)	Heptachlor (76-44-8)	1,1,1,2-Tetrachloroethane (630-20-6)
2-Chloronaphthalene (91-58-7)	Heptachlor epoxide (1024-57-3)	1,1,2,2-Tetrachloroethane (79-34-5)
2-Chlorophenol (95-57-8)	Hexachlorobenzene (118-74-1)	Tetrachloroethylene; Tetrachloroethene; Perchloroethylene (127-18-4)
4-Chlorophenyl phenyl ether (7005-72-3)	Hexachlorobutadiene (87-68-3)	2,3,4,6-Tetrachlorophenol (58-90-2)
Chloroprene (126-99-8)	Hexachlorocyclopentadiene (77-47-4)	Toluene (108-88-3)
Chrysene (218-01-9)	Hexachloroethane (67-72-1)	o-Toluidine (95-53-4)
m-Cresol, 3-methylphenol (108-39-4)	Hexachloropropene (1888-71-7)	Toxaphene <sup>28</sup>
o-Cresol, 2-methylphenol (95-48-7)	2-Hexanone, Methyl butyl ketone (591-78-6)	1,2,4-Trichlorobenzene (120-82-1)
p-Cresol; 4-methylphenol (106-44-5)	Indeno(1,2,3-cd)pyrene (193-39-5)	1,1,1-Trichloroethane, Methylchloroform (71-55-6)
2,4-D, 2,4-Dichlorophen- oxyacetic acid (94-75-7)	Isobutyl alcohol (78-83-1)	1,1,2-Trichloroethane (79-00-5)
4,4 <sup>21</sup> -DDD (72-54-8)	Isodrin (465-73-6)	Trichloroethylene, Trichloroethene (79-01-6)
4,4 <sup>21</sup> -DDE (72-55-9)	Isophorone (78-59-1)	Trichlorofluoromethane, R-11 (75-69-4)
4,4 <sup>21</sup> -DDT (50-29-3)	Isosafrole (120-58-1)	2,4,5-Trichlorophenol (95-95-4)
Diallate (2303-16-4)	Kepone (143-50-0)	2,4,6-Trichlorophenol (88-06-2)
Dibenz[a,h]anthracene (53-70-3)	Methacrylonitrile (126-98-7)	1,2,3-Trichloropropane (96-18-4)
Dibenzofuran (132-64-9)	Methapyrilene (91-80-5)	0,0,0-Triethyl phosphorothioate (126-68-1)
Dibromochloromethane; Chlorodibromomethane (124-48-1)	Methoxychlor (72-43-5)	sym-Trinitrobenzene (99-35-4)
1,2-Dibromo-3-chloro- propane; DBCP (96-12-8)	Methyl bromide, Bromomethane (74-83-9)	Vinyl acetate (108-05-4)
1,2-Dibromoethane, Ethylene dibromide; EDB (106-93-4)	Methyl chloride, Chloromethane (74-87-3)	Vinyl chloride; Chloroethene (75-01-4)
Di-n-butyl phthalate (84-74-2)	3-Methylcholanthrene (56-49-5)	Xylene (total)
o-Dichlorobenzene; 1,2-Dichlorobenzene (95-50-1)	Methyl ethyl ketone, MEK, 2-Butanone (78-93-3)	Per- and polyfluoroalkyl substances <sup>29</sup>
m-Dichlorobenzene; 1,3-Dichlorobenzene (541-73-1)	Methyl iodide, Iodomethane (74-88-4)	1,4-Dioxane (123-91-1)
p-Dichlorobenzene; 1,4-dichlorobenzene (106-46-7)	Methyl methacrylate (80-62-6)	
3,3 <sup>21</sup> -Dichlorobenzidine (91-94-1)	Methyl methanesulfonate (66-27-3)	
trans-1,4-Dichloro- 2-butene (110-57-6)	2-Methylnaphthalene (91-57-6)	

Dichlorodifluoromethane, CFC 12 (75-71-8)	Methyl parathion; Parathion methyl (298-00-0)
1,1-Dichloroethane; Ethylidene chloride (75-34-3)	4-Methyl-2-pentanone, Methyl isobutyl ketone (108-10-1)
1,2-Dichloroethane; Ethylene dichloride (107-06-2)	Methylene bromide; Dibromomethane (74-95-3)
1,1-Dichloroethylene, 1,1-Dichloroethene; Vinylidene chloride (75-35-4)	Methylene chloride, Dichloromethane (75-09-2)

**(i) Leachate management plan.**

The leachate management plan must include:

- (1) a description of how the landfill will be constructed, operated, and closed in a manner that minimizes the generation of leachate, except in those cases where the department has approved the recirculation of leachate for waste mass stabilization enhancement, and how the migration of leachate into surface water or groundwater will be prevented;
- (2) a description of operational methods to minimize the occurrence of perched leachate trapped above the leachate collection and removal system and surface seeps of leachate from above-grade landfill operations;
- (3) a schedule for biennial video inspection and annual maintenance of the primary and secondary leachate collection and removal system;
- (4) a schedule for the monitoring and recording of the secondary leachate collection and removal system flow data to determine the presence, quantity, nature and significance of any liquid detected;
- (5) a discussion of the specific design and operational features related to the system, including leachate monitoring and sampling, locations of all leachate sampling points, alarm systems and maintenance, and any required back up equipment; and
- (6) if leachate recirculation is proposed, the leachate management plan must include
  - (i) a supporting geotechnical analysis evaluating the effect of leachate recirculation on the structural integrity and stability of the landfill's liner system, leachate collection and removal system, and waste mass;
  - (ii) a description of how increased landfill gas emissions and associated odors will be controlled;
  - (iii) a description of the methods and rate of leachate recirculation and addition;
  - (iv) procedures for recording the date and volume of recirculated leachate;
  - (v) a description of the operation, which addresses:
    - (a) the use of permeable operating cover or alternative operating cover to facilitate leachate distribution throughout the waste mass, and
    - (b) operational controls such as monitoring of surface seeps, liner system performance and excessive leachate head buildup, prevention of subsurface fires, odor control, and instruction for cessation of leachate recirculation and remediation of these conditions.

**(j) Odor control plan.**

The odor control plan must include:

- (1) identification of all potential sources for odors and a description of the operational procedures and strategies to be followed to effectively control odors at the facility;
- (2) procedures to be taken in the event of proposed waste volume increases or changes in waste characterization that may increase landfill gas emissions or odors;
- (3) identification of the landfill personnel who would be responsible for implementation of the odor control plan; and
- (4) operational and design-related recommendations that can be implemented upon detection of odor control problems, including impervious membranes and interim covers in conjunction with other landfill gas control methods. The odor control plan may include but not be limited to, gas control systems that are appropriately connected to the landfill liner system's primary leachate collection and removal system (including the drainage area on the landfill's side slopes), use of a horizontal gas collection lines, which may include rejection or mitigation of odiferous wastes that are determined to be contributing to off-site odors.

**(k) Gas monitoring and emission control plan.**

The gas monitoring and emission control plan must include:

- (1) a description of the day-to-day operation of the landfill gas management system with respect to operation of odor and emission controls;

(2) a description of any air quality monitoring, including monitoring for fugitive landfill odor and air emissions; and

(3) for a landfill with an appurtenant landfill gas-to-energy facility or other landfill gas recovery facility, a discussion of how the landfill's odor and air emission controls are integrated with a recovery facility.

**(l) Winter and inclement weather operation plan.**

A description of how winter and inclement weather operations will be conducted, including identification of the specific actions to be taken to prevent frost action on the liner system in places where waste will not be placed within one year of construction certification approval.

**(m) Residential drop-off operation plan.**

A description of the operation of a residential drop-off area, if applicable, for non-commercial vehicles to unload waste and recyclables at an area other than the landfill working face.

**(n) A radioactive waste detection plan.**

The radioactive waste detection plan must include procedures for detecting radioactive material; operation and maintenance documents for radiation detectors which address proper equipment placement for effective operation and include setting of investigation alarm setpoint settings and calibration methods; and response procedures to be implemented if radioactive waste is detected.

**(o) Emergency response plan.**

An emergency response plan must include a description of, at a minimum, the actions to be taken in response to:

- (1) uncontrolled explosive landfill gases detected on-site or beyond the property boundary;
- (2) unexpected events during the construction and operation of the landfill gas management system, including the equipment to be utilized to maintain proper landfill gas venting and control when normal operations cease; and
- (3) unexpected events during the subsequent construction and/or daily operation of the landfill's leachate collection and removal system.

**(p) Conceptual closure, post-closure care, custodial care, and end use plan.**

The conceptual closure, post-closure care, custodial care, and end use plan must include:

- (1) a site plan that shows proposed final contours, property lines, storm water drainage system, streams and water courses, roads, structures and, if applicable, the groundwater and leachate treatment system, air pollution control system and any active landfill gas collection system;
- (2) typical details of final cover system components and facility structures;
- (3) a description of how the sequential closure of areas of the landfill is expected to progress in concert with the fill progression schedule, including effects of landfill reclamation activities if proposed;
- (4) an estimate of the greatest number of landfill cells which, at any given point during the lifetime of the facility, will have received waste but not undergone final closure;
- (5) an estimate of the maximum volume of waste and alternative operating cover that will be contained within the landfill;
- (6) sufficient information upon which to estimate closure costs and post-closure and custodial care monitoring and maintenance costs. This information must be based upon the requirements of Subpart 363-9 of this Part, including a rolling 30-year post-closure care period, and must include estimates of:
  - (i) quantities and costs for each component of the final cover system, including related construction costs;
  - (ii) the anticipated length of the post-closure care period based on the types of wastes disposed and the criteria provided in section 363-9.6(a) of this Part;
  - (iii) post-closure operational, monitoring and maintenance costs including costs to replace system components based on predicted service life; and
  - (iv) custodial care monitoring and maintenance costs including costs to replace system components based on predicted service life; and
- (7) a conceptual end use for the site, if proposed.

**Footnotes**

- 1 This list contains parameters for which possible analytical procedures are provided in: *Test Methods for Evaluating Solid Waste, Physical/Chemical Methods*, EPA Publication SW-846 (Third Edition, (November 1986), as amended by Updates I

- (July 1992), II (September 1994), IIA (August 1993), IIB (January 1995), III (December 1996), IIIA (April 1998), document number 955-001-00000-1), incorporated by reference in section 360.3 of this Title. *Methods for Chemical Analysis of Water and Wastes*, USEPA-600/4-79-020, March, 1983, incorporated by reference in section 360.3 of this Title.
- 2 Common names are those widely used in government regulations, scientific publications, and commerce; synonyms exist for many chemicals. "Total" indicates all species in the groundwater that contain this element.
- 3 Any floaters or sinkers found must be analyzed separately for baseline parameters.
- 4 Surface water only.
- 5 Any unusual conditions (colors, odors, surface sheens, etc.) noticed during well development, purging, or sampling must be reported.
- 6 This list contains parameters for which possible analytical procedures are provided in: *Test Methods for Evaluating Solid Waste, Physical/Chemical Methods*, EPA Publication SW-846 (Third Edition, (November 1986), as amended by Updates I (July 1992), II (September 1994), IIA (August 1993), IIB (January 1995), III (December 1996), IIIA (April 1998), document number 955-001-00000-1), incorporated by reference in section 360.3 of this Title. *Methods for Chemical Analysis of Water and Wastes*, USEPA-600/4-79-020, March, 1983, incorporated by reference in section 360.3 of this Title.
- 7 Common names are those widely used in government regulations, scientific publications, and commerce; synonyms exist for many chemicals. "Total" indicates all species in the groundwater that contain this element.
- 8 Any floaters or sinkers found must be analyzed separately for baseline parameters.
- 9 Surface water only.
- 10 Any unusual conditions (colors, odors, surface sheens, etc.) noticed during well development, purging, or sampling must be reported.
- 11 The department may waive the requirement to analyze hexavalent chromium provided that total and hexavalent and trivalent chromium values do not exceed 0.05 mg/l.
- 12 This list contains parameters for which possible analytical procedures are provided in: *Test Methods for Evaluating Solid Waste, Physical/Chemical Methods*, EPA Publication SW-846 (Third Edition, (November 1986), as amended by Updates I (July 1992), II (September 1994), IIA (August 1993), IIB (January 1995), III (December 1996), and IIIA (April 1998) document number 955-001-00000-1), incorporated by reference in section 360.3 of this Title. *Methods for Chemical Analysis of Water and Wastes*, USEPA-600/4-79-020, March, 1983, incorporated by reference in 360.3 of this Title.
- 13 Common names are those widely used in government regulations, scientific publications, and commerce; synonyms exist for many chemicals.
- 14 This list contains parameters for which possible analytical procedures are provided in: *Test Methods for Evaluating Solid Waste, Physical/Chemical Methods*, EPA Publication SW-846 (Third Edition, (November 1986), as amended by Updates I (July 1992), II (September 1994), IIA (August 1993), IIB (January 1995), III (December 1996), and IIIA (April 1998) document number 955-001-00000-1), incorporated by reference in section 360.3 of this Title. *Methods for Chemical Analysis of Water and Wastes*, USEPA-600/4-79-020, March 1983, incorporated by reference in 360.3 of this Title. *Prescribed Procedures for Measurement of Radioactivity in Drinking Water*, USEPA-600/4-80-032, August 1980, incorporated by reference in section 360.3 of this Title.
- 15 Common names are those widely used in government regulations, scientific publications, and commerce; synonyms exist for many chemicals. "Total" indicates all species in the groundwater that contain this element.
- 16 Two sets of samples must be collected: one filtered and one unfiltered. Filtered samples must be filtered using a 0.45 micron filter via standard techniques.
- 17 Any floaters or sinkers found must be analyzed separately for baseline parameters.
- 18 Surface water only.
- 19 Any unusual conditions (colors, odors, surface sheens, etc.) noticed during well development, purging, or sampling must be reported.
- 20 The department may waive the requirement to analyze hexavalent chromium provided that total and hexavalent and trivalent chromium values do not exceed 0.05 mg/l.
- 21 This list contains parameters for which possible analytical procedures are provided in: *Test Methods for Evaluating Solid Waste, Physical/Chemical Methods*, EPA Publication SW-846 (Third Edition, (November 1986), as amended by Updates I (July 1992), II (September 1994), IIA (August 1993), IIB (January 1995), III (December 1996), and IIIA (April 1998) document number 955-001-00000-1), incorporated by reference in section 360.3 of this Title. *Methods for Chemical Analysis of Water and Wastes*, USEPA-600/4-79-020, March 1983, incorporated by reference in section 360.3 of this Title.

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Common names are those widely used in government regulations, scientific publications, and commerce; synonyms exist for many chemicals.

- 23 This substance is often called Bis(2-chloroisopropyl) ether, the name Chemical Abstracts Service applies to its noncommercial isomer, Propane, 2,2"-oxybis[2]-chloro- (CAS RN 39638-32-9).
- 24 Polychlorinated biphenyls (1336-36-3): This category contains congener chemicals, including constituents of Aroclor 1016 (12674-11-2), Aroclor 1221 (11104-28-2), Aroclor 1232 (11097-69-1), and Aroclor 1260 (11096-82-5).
- 25 Polychlorinated dibenzo-p-dioxins: This category contains congener chemicals, including tetrachlorodibenzo-p-dioxins, pentachlorodibenzo-p-dioxins, and hexachlorodibenzo-p-dioxins.
- 26 Polychlorinated dibenzofurans: This category includes congener chemicals, including tetrachlorodibenzofurans, pentachlorodibenzofurans, and hexachlorodibenzofurans.
- 27 Chlordane: This entry includes alpha-chlordane (5103-71-9), beta-chlordane (5103-74-2), gamma-chlordane (5566-34-7), and constituents of chlordane (57-74-9; 12789-03-6).
- 28 Toxaphene: This entry includes congener chemicals contained in technical toxaphene (CAS RN 8001-35-2), *i.e.*, chlorinated camphene.
- 29 Per- and polyfluoroalkyl substances (PFAS): This category contains congener chemicals, including but not limited to perfluorooctanoic acid, perfluorooctanesulfonic acid, perfluorononanoic acid, perfluorohexanesulfonic acid, perfluoroheptanoic acid, perfluorobutanesulfonic acid.

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