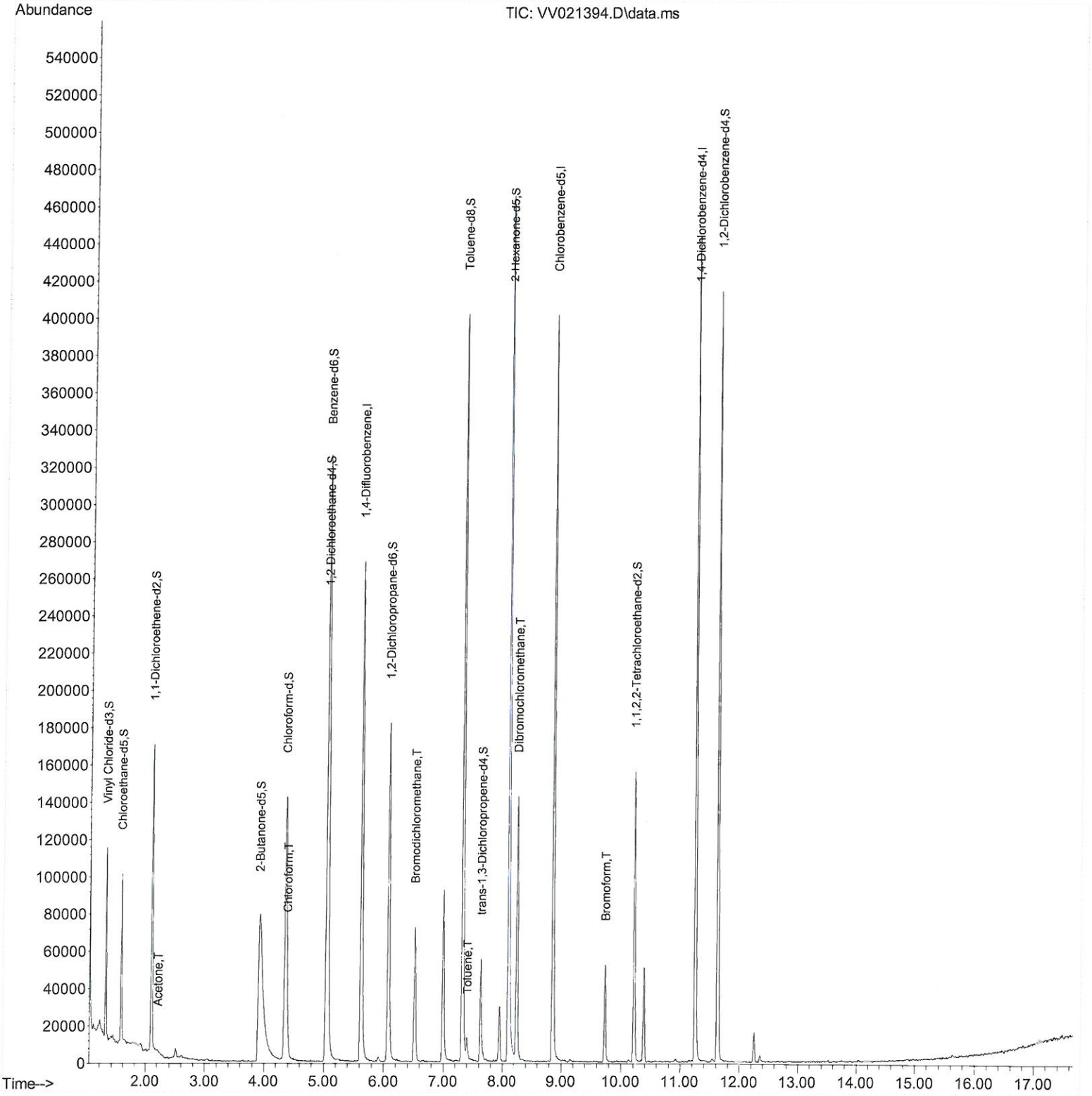
Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV052521\
Data File : VV021394.D
Acq On : 25 May 2021 17:37
Operator : SY/MD
Sample : M2496-13
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 16 Sample Multiplier: 1

Quant Time: May 26 07:54:01 2021
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR052521WMA.M
Quant Title : TRACE VOA SFAM1.0
QLast Update : Wed May 26 07:52:48 2021
Response via : Initial Calibration

Instrument :
MSVOA_V
ClientSampleId :
GB444

Manual Integrations APPROVED
MMDadoda
5/26/2021 3:13:19 PM



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV052521\
Data File : VV021394.D
Acq On : 25 May 2021 17:37
Operator : SY/MD
Sample : M2496-13
Misc : 25.0mL/MSVOA_V/WATER
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Quant Title : TRACE VOA SFAM1.0
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MMDadoda
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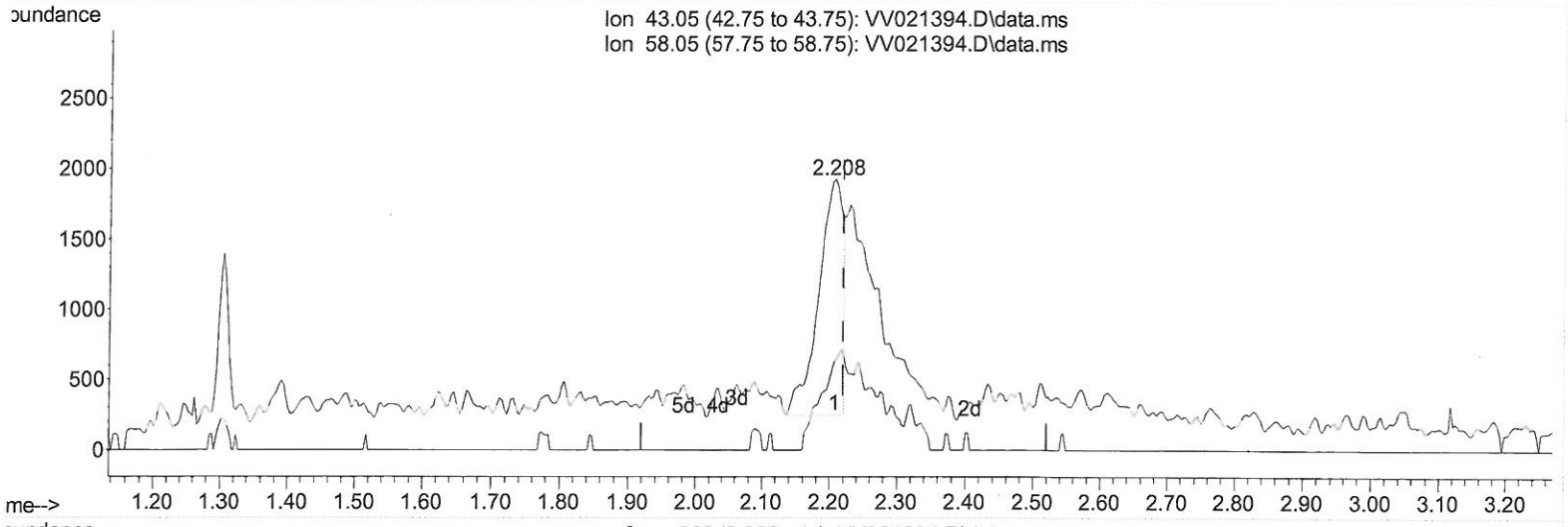

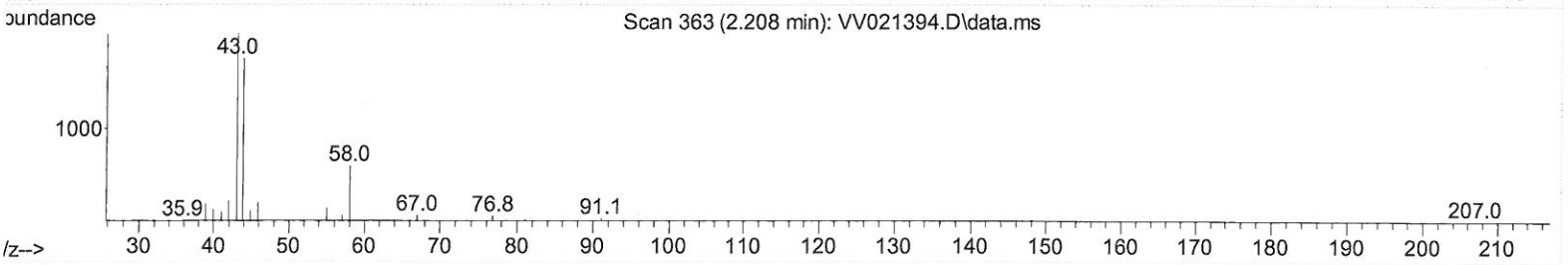

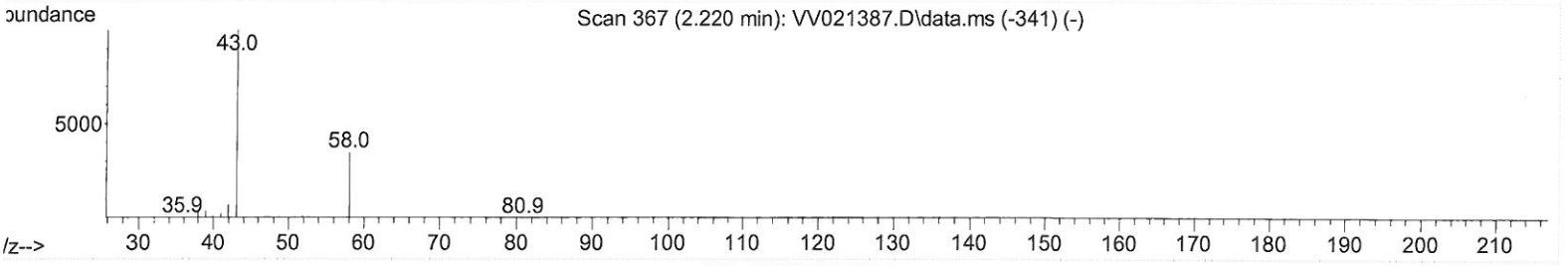

TIC: VV021394.D\data.ms

(13) Acetone (T)

2.208min (-0.013) 1.93 ug/L

response 4197

Ion	Exp%	Act%
43.05	100.00	100.00
58.05	9.30	47.84#
0.00	0.00	0.00
0.00	0.00	0.00

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV052521\
Data File : VV021394.D
Acq On : 25 May 2021 17:37
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Misc : 25.0mL/MSVOA_V/WATER
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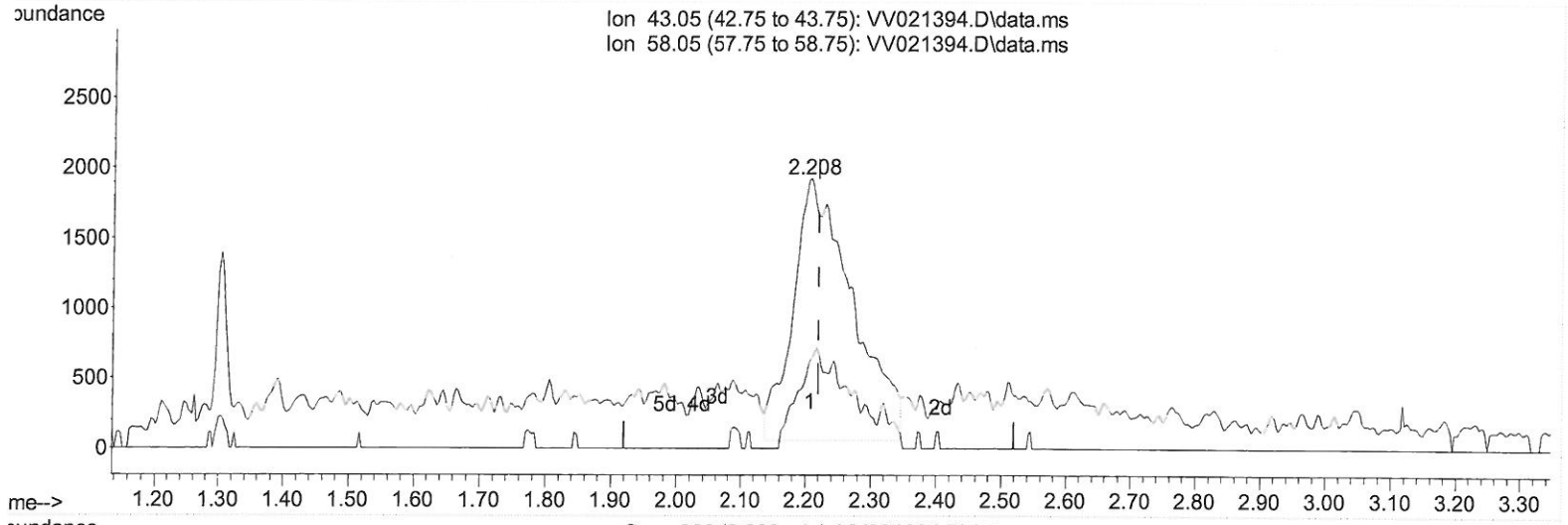
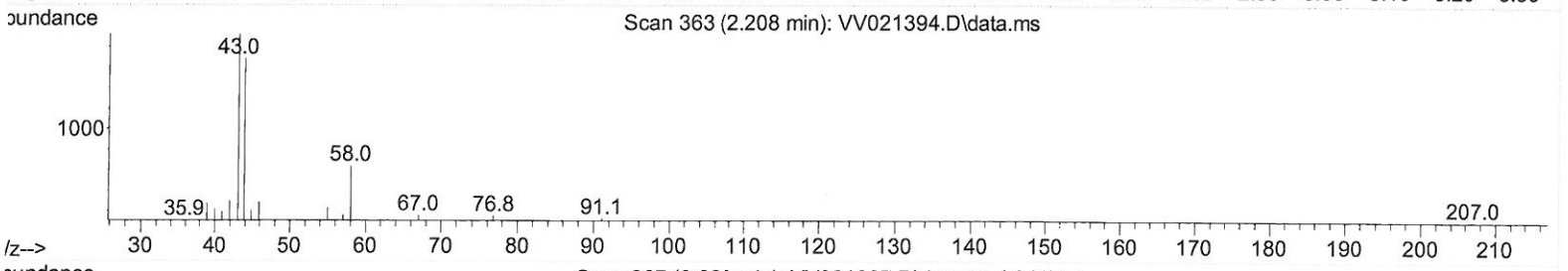
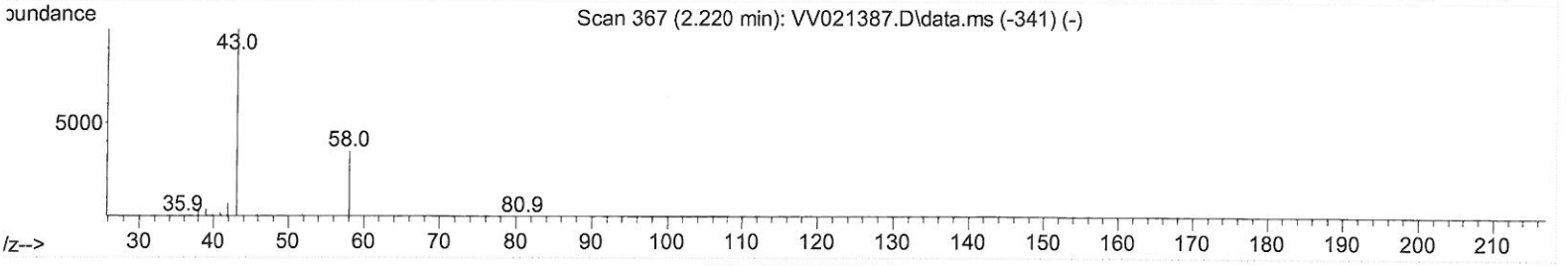
Quant Time: May 26 07:54:01 2021
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR052521WMA.M
Quant Title : TRACE VOA SFAM1.0
QLast Update : Wed May 26 07:52:48 2021
Response via : Initial Calibration

Instrument :
MSVOA_V
ClientSampleId :
GB444

Manual Integrations APPROVED
MMDadoda
5/26/2021 3:13:19 PM







TIC: VV021394.D\data.ms

(13) Acetone (T)

2.208min (-0.013) 5.44 ug/L m

response 11812

Ion	Exp%	Act%
43.05	100.00	100.00
58.05	9.30	17.00
0.00	0.00	0.00
0.00	0.00	0.00

7 MD 5/28/21

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV052521\
Data File : VV021394.D
Acq On : 25 May 2021 17:37
Operator : SY/MD
Sample : M2496-13
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 16 Sample Multiplier: 1

Quant Time: May 26 07:54:01 2021
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR052521WMA.M
Quant Title : TRACE VOA SFAM1.0
QLast Update : Wed May 26 07:52:48 2021
Response via : Initial Calibration

Instrument : MSVOA_V
ClientSampleId : GB444

Manual Integrations APPROVED
MMDadoda
5/26/2021 3:13:19 PM

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Difluorobenzene	5.619	114	253083	5.00	ug/L	0.00
28) Chlorobenzene-d5	8.854	117	233638	5.00	ug/L	0.00
58) 1,4-Dichlorobenzene-d4	11.249	152	127336	5.00	ug/L	0.00
System Monitoring Compounds						
4) Vinyl Chloride-d3	1.304	65	63025	5.89	ug/L	0.00
Spiked Amount 5.000	Range 40 - 130		Recovery = 117.80%			
7) Chloroethane-d5	1.564	69	56891	5.57	ug/L	0.00
Spiked Amount 5.000	Range 65 - 130		Recovery = 111.40%			
11) 1,1-Dichloroethene-d2	2.105	63	82899	4.08	ug/L	0.00
Spiked Amount 5.000	Range 60 - 125		Recovery = 81.60%			
20) 2-Butanone-d5	3.915	46	207978	54.03	ug/L	0.00
Spiked Amount 50.000	Range 40 - 130		Recovery = 108.06%			
24) Chloroform-d	4.352	84	150464	5.07	ug/L	0.00
Spiked Amount 5.000	Range 70 - 125		Recovery = 101.40%			
26) 1,2-Dichloroethane-d4	5.037	65	72223	5.29	ug/L	0.00
Spiked Amount 5.000	Range 70 - 130		Recovery = 105.80%			
32) Benzene-d6	5.053	84	301591	5.37	ug/L	0.00
Spiked Amount 5.000	Range 70 - 125		Recovery = 107.40%			
36) 1,2-Dichloropropane-d6	6.072	67	95827	5.26	ug/L	0.00
Spiked Amount 5.000	Range 60 - 140		Recovery = 105.20%			
41) Toluene-d8	7.320	98	272838	5.29	ug/L	0.00
Spiked Amount 5.000	Range 70 - 130		Recovery = 105.80%			
43) trans-1,3-Dichloroprop...	7.629	79	35202	4.64	ug/L	0.00
Spiked Amount 5.000	Range 55 - 130		Recovery = 92.80%			
46) 2-Hexanone-d5	8.092	63	183451	51.35	ug/L	0.00
Spiked Amount 50.000	Range 45 - 130		Recovery = 102.70%			
56) 1,1,2,2-Tetrachloroeth...	10.217	84	75075	5.11	ug/L	0.00
Spiked Amount 5.000	Range 65 - 120		Recovery = 102.20%			
66) 1,2-Dichlorobenzene-d4	11.625	152	115526	5.26	ug/L	0.00
Spiked Amount 5.000	Range 80 - 120		Recovery = 105.20%			
Target Compounds						Qvalue
13) Acetone	2.208	43	11812m	5.44	ug/L	
25) Chloroform	4.381	83	19890	0.70	ug/L	98
38) Bromodichloromethane	6.516	83	50843	2.46	ug/L	95
42) Toluene	7.400	91	8520	0.13	ug/L	97
49) Dibromochloromethane	8.249	129	76926	5.07	ug/L	100
59) Bromoform	9.735	173	24963	2.71	ug/L	98

(#) = qualifier out of range (m) = manual integration (+) = signals summed

> MD 5/28/21