

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN111921\
 Data File : VN069599.D
 Acq On : 19 Nov 2021 21:12
 Operator : JC/MD
 Sample : M4764-03
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 MW-3R

Quant Time: Nov 22 05:10:06 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N110921W.M
 Quant Title : SW846 8260
 QLast Update : Tue Nov 09 18:18:22 2021
 Response via : Initial Calibration

Compound	R.T. QIon		Response	Conc	Units	Dev(Min)

Internal Standards						
1) Pentafluorobenzene	8.088	168	1400418	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.968	114	2443850	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.746	117	2468980	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.675	152	926713	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.440	65	1043472	51.389	ug/l	0.00
Spiked Amount	50.000	Range 61 - 141	Recovery	= 102.780%		
35) Dibromofluoromethane	8.024	113	718615	49.028	ug/l	0.00
Spiked Amount	50.000	Range 69 - 133	Recovery	= 98.060%		
50) Toluene-d8	10.443	98	3132546	54.341	ug/l	0.00
Spiked Amount	50.000	Range 65 - 126	Recovery	= 108.680%		
62) 4-Bromofluorobenzene	12.733	95	1132216	53.464	ug/l	0.00
Spiked Amount	50.000	Range 58 - 135	Recovery	= 106.920%		
Target Compounds						
						Qvalue
7) Trichlorofluoromethane	3.381	101	39154	1.458	ug/l	99
44) Trichloroethene	9.217	130	129913	8.552	ug/l	94
64) Tetrachloroethene	10.977	164	2637072	176.890	ug/l	95

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN111921\
Data File : VN069599.D
Acq On : 19 Nov 2021 21:12
Operator : JC/MD
Sample : M4764-03
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-3R

Quant Time: Nov 22 05:10:06 2021
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N110921W.M
Quant Title : SW846 8260
QLast Update : Tue Nov 09 18:18:22 2021
Response via : Initial Calibration

