

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN111621\
 Data File : VN069518.D
 Acq On : 16 Nov 2021 19:23
 Operator : JC/MD
 Sample : M4714-03
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 AEI-1

Quant Time: Nov 17 02:01:57 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	8.086	168	1419177	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.968	114	2477114	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.747	117	2447865	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.678	152	919963	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.440	65	1034546	50.276	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	= 100.560%	
35) Dibromofluoromethane	8.024	113	718634	48.371	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	= 96.740%	
50) Toluene-d8	10.443	98	3144165	53.810	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	= 107.620%	
62) 4-Bromofluorobenzene	12.734	95	1112879	51.845	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	= 103.700%	
Target Compounds						Qvalue
3) Chloromethane	2.306	50	3568	0.203	ug/l	95
27) cis-1,2-Dichloroethene	7.327	96	17366	1.090	ug/l	100
30) Chloroform	7.823	83	17541	0.613	ug/l	94
44) Trichloroethene	9.217	130	11164	0.725	ug/l	90
64) Tetrachloroethene	10.977	164	29400	1.989	ug/l	94

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

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