

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN102721\
 Data File : VN069225.D
 Acq On : 27 Oct 2021 18:03
 Operator : JC/MD
 Sample : M4338-06
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 MW-2D(ONSITE)

Quant Time: Oct 28 05:22:24 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N101221W.M
 Quant Title : SW846 8260
 QLast Update : Wed Oct 13 05:44:40 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.088	168	323229	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.971	114	571829	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.747	117	543691	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.678	152	212207	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.440	65	215178	50.492	ug/l	0.00
Spiked Amount 50.000	Range 61 - 141		Recovery = 100.980%			
35) Dibromofluoromethane	8.027	113	165524	49.752	ug/l	0.00
Spiked Amount 50.000	Range 69 - 133		Recovery = 99.500%			
50) Toluene-d8	10.443	98	709493	53.299	ug/l	0.00
Spiked Amount 50.000	Range 65 - 126		Recovery = 106.600%			
62) 4-Bromofluorobenzene	12.734	95	258908	51.551	ug/l	0.00
Spiked Amount 50.000	Range 58 - 135		Recovery = 103.100%			
Target Compounds						
30) Chloroform	7.823	83	40515	6.077	ug/l	96
38) Carbon Tetrachloride	8.212	117	70222	13.091	ug/l	98
44) Trichloroethene	9.218	130	11173	3.002	ug/l	95
64) Tetrachloroethene	10.985	164	12605408	3599.083	ug/l	96
90) Hexachloroethane	14.211	117	232138	77.677	ug/l	99

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

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