

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN102721\
 Data File : VN069207.D
 Acq On : 27 Oct 2021 1:10
 Operator : JC/MD
 Sample : M4337-21
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 DDC-2-PD

Quant Time: Oct 27 09:26:18 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	325685	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.971	114	568643	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.746	117	552088	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.677	152	205456	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.440	65	213570	49.736	ug/l	0.00
Spiked Amount 50.000	Range 61 - 141		Recovery =	99.480%		
35) Dibromofluoromethane	8.027	113	164963	49.862	ug/l	0.00
Spiked Amount 50.000	Range 69 - 133		Recovery =	99.720%		
50) Toluene-d8	10.446	98	712118	53.795	ug/l	0.00
Spiked Amount 50.000	Range 65 - 126		Recovery =	107.600%		
62) 4-Bromofluorobenzene	12.733	95	257350	51.528	ug/l	0.00
Spiked Amount 50.000	Range 58 - 135		Recovery =	103.060%		
Target Compounds						
					Qvalue	
16) Acetone	4.307	43	5630	3.513	ug/l #	89
30) Chloroform	7.817	83	4365	0.650	ug/l #	82
38) Carbon Tetrachloride	8.209	117	6356	1.192	ug/l	95
44) Trichloroethene	9.220	130	3032	0.819	ug/l	97
64) Tetrachloroethene	10.982	164	2430100	683.287	ug/l	98
90) Hexachloroethane	14.214	117	48706	16.833	ug/l	99

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

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