

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN102921\
 Data File : VN069296.D
 Acq On : 29 Oct 2021 13:04
 Operator : JC/MD
 Sample : M4338-07
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 8 Sample Multiplier: 1

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

Quant Time: Oct 30 05:40:37 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N102821W.M
 Quant Title : SW846 8260
 QLast Update : Thu Oct 28 16:09:15 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.088	168	317907	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.971	114	556718	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.746	117	544090	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.677	152	205063	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.440	65	213851	46.600	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	=	93.200%
35) Dibromofluoromethane	8.024	113	161300	46.479	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	=	92.960%
50) Toluene-d8	10.443	98	699260	50.395	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	=	100.780%
62) 4-Bromofluorobenzene	12.733	95	251078	47.952	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	=	95.900%
Target Compounds						
					Qvalue	
16) Acetone	4.299	43	3389	1.733	ug/l	99
64) Tetrachloroethene	10.982	164	2840	0.817	ug/l	91

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

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