

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN102921\
 Data File : VN069298.D
 Acq On : 29 Oct 2021 13:55
 Operator : JC/MD
 Sample : M4338-15
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 MW-1D(ONSITE)

Manual Integrations
 APPROVED

Reviewed By :Mahesh Dadoda 11/01/2021
 Supervised By :Semsettin Yesilyurt 11/01/2021

Quant Time: Oct 30 05:41:16 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	312289	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.968	114	551989	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.747	117	531519	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.678	152	199753	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.440	65	211862	46.997	ug/l	0.00
Spiked Amount 50.000	Range 61 - 141		Recovery =	94.000%		
35) Dibromofluoromethane	8.024	113	159984	46.495	ug/l	0.00
Spiked Amount 50.000	Range 69 - 133		Recovery =	92.980%		
50) Toluene-d8	10.443	98	685977	49.861	ug/l	0.00
Spiked Amount 50.000	Range 65 - 126		Recovery =	99.720%		
62) 4-Bromofluorobenzene	12.733	95	244222	47.042	ug/l	0.00
Spiked Amount 50.000	Range 58 - 135		Recovery =	94.080%		
Target Compounds					Qvalue	
16) Acetone	4.291	43	3735m	1.945	ug/l	
44) Trichloroethene	9.215	130	2534	0.703	ug/l	92
64) Tetrachloroethene	10.977	164	3571	1.052	ug/l	95

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

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