

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN102921\
 Data File : VN069301.D
 Acq On : 29 Oct 2021 15:13
 Operator : JC/MD
 Sample : M4338-17
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 MW-14S

Manual Integrations
 APPROVED

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

Quant Time: Oct 30 05:42:14 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	292791	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.971	114	520004	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.747	117	506225	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.678	152	189482	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.440	65	199526	47.208	ug/l	0.00
Spiked Amount 50.000	Range 61 - 141		Recovery =	94.420%		
35) Dibromofluoromethane	8.024	113	149117	46.002	ug/l	0.00
Spiked Amount 50.000	Range 69 - 133		Recovery =	92.000%		
50) Toluene-d8	10.443	98	650738	50.209	ug/l	0.00
Spiked Amount 50.000	Range 65 - 126		Recovery =	100.420%		
62) 4-Bromofluorobenzene	12.734	95	237693	48.601	ug/l	0.00
Spiked Amount 50.000	Range 58 - 135		Recovery =	97.200%		
Target Compounds						
16) Acetone	4.296	43	4103	2.278	ug/l	96
30) Chloroform	7.817	83	3039	0.502	ug/l	84
44) Trichloroethene	9.212	130	1033m	0.304	ug/l	
64) Tetrachloroethene	10.980	164	221306	68.442	ug/l	99
90) Hexachloroethane	14.211	117	3397	3.323	ug/l	99

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

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