

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN102921\
 Data File : VN069314.D
 Acq On : 29 Oct 2021 20:47
 Operator : JC/MD
 Sample : M4390-01
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 FRAC-COMP

Quant Time: Oct 30 05:47:07 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

Manual Integrations
 APPROVED

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

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

Internal Standards						
1) Pentafluorobenzene	8.088	168	310152	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.971	114	553350	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.747	117	533602	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.678	152	199423	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.440	65	213748	47.742	ug/l	0.00
Spiked Amount 50.000	Range 61 - 141		Recovery =	95.480%		
35) Dibromofluoromethane	8.024	113	158454	45.937	ug/l	0.00
Spiked Amount 50.000	Range 69 - 133		Recovery =	91.880%		
50) Toluene-d8	10.443	98	690072	50.035	ug/l	0.00
Spiked Amount 50.000	Range 65 - 126		Recovery =	100.080%		
62) 4-Bromofluorobenzene	12.733	95	247285	47.515	ug/l	0.00
Spiked Amount 50.000	Range 58 - 135		Recovery =	95.040%		
Target Compounds						
16) Acetone	4.299	43	4711m	2.470	ug/l	
30) Chloroform	7.820	83	40342	6.292	ug/l	96
47) Bromodichloromethane	9.767	83	2917	1.619	ug/l #	96

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

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