

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN041522\
 Data File : VN072022.D
 Acq On : 15 Apr 2022 23:40
 Operator : JC\MD
 Sample : PB144101ZHE#10
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB144101ZHE#10

Manual Integrations
 APPROVED

Reviewed By :Mahesh Dadoda 04/18/2022
 Supervised By :Semsettin Yesilyurt 04/18/2022

Quant Time: Apr 16 01:07:17 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N033022W.M
 Quant Title : SW846 8260
 QLast Update : Thu Mar 31 01:57:54 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	7.710	168	607640m	50.000	ug/l	-0.37
34) 1,4-Difluorobenzene	8.710	114	1051548	50.000	ug/l	-0.25
63) Chlorobenzene-d5	11.686	117	1227760m	50.000	ug/l	-0.05
72) 1,4-Dichlorobenzene-d4	13.657	152	469391	50.000	ug/l	-0.01
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.098	65	332666	47.956	ug/l	-0.34
Spiked Amount 50.000	Range 61 - 141		Recovery =	95.920%		
35) Dibromofluoromethane	7.616	113	290513	46.950	ug/l	-0.40
Spiked Amount 50.000	Range 69 - 133		Recovery =	93.900%		
50) Toluene-d8	10.316	98	1371699	54.190	ug/l	-0.12
Spiked Amount 50.000	Range 65 - 126		Recovery =	108.380%		
62) 4-Bromofluorobenzene	12.698	95	510972m	62.416	ug/l	-0.03
Spiked Amount 50.000	Range 58 - 135		Recovery =	124.840%		
Target Compounds					Qvalue	
16) Acetone	3.275	43	17019m	5.741	ug/l	
43) Isopropyl Acetate	8.263	43	53701	3.414	ug/l #	92
71) Bromoform	12.698	173	3872	0.573	ug/l #	1

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

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