

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN062123\
 Data File : VN078322.D
 Acq On : 21 Jun 2023 15:10
 Operator : JC\MD
 Sample : 03359-01
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 OWBR-03-128-148-062023

Quant Time: Jun 22 04:50:01 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N060623W.M
 Quant Title : SW846 8260
 QLast Update : Wed Jun 07 05:05:00 2023
 Response via : Initial Calibration

Manual Integrations
 APPROVED

Reviewed By : John Carlone 06/22/2023
 Supervised By : Mahesh Dadoda 06/22/2023

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

Internal Standards						
1) Pentafluorobenzene	8.235	168	417660	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.112	114	782304	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	707497	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.800	152	217043	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.588	65	298401	53.586	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery = 107.180%			
35) Dibromofluoromethane	8.177	113	274437	55.482	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery = 110.960%			
50) Toluene-d8	10.571	98	925814	48.551	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery = 97.100%			
62) 4-Bromofluorobenzene	12.859	95	294393	49.661	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery = 99.320%			
Target Compounds						
17) Carbon Disulfide	4.730	76	9817m	0.836	ug/l	
40) Benzene	8.612	78	81212	3.834	ug/l	100
67) Ethyl Benzene	11.965	91	12157	0.465	ug/l	96

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

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