

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN072424\
 Data File : VN083028.D
 Acq On : 24 Jul 2024 14:53
 Operator : JC\MD
 Sample : P3338-03
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 VPB196-HYD-20240718

Quant Time: Jul 25 02:39:12 2024
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N070824W.M
 Quant Title : SW846 8260
 QLast Update : Mon Jul 08 13:46:04 2024
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.223	168	149005	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	289318	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.864	117	295637	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	124529	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.582	65	120400	57.642	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery = 115.280%			
35) Dibromofluoromethane	8.171	113	101314	54.573	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery = 109.140%			
50) Toluene-d8	10.565	98	341068	49.741	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery = 99.480%			
62) 4-Bromofluorobenzene	12.853	95	114309	46.410	ug/l	0.00
Spiked Amount 50.000	Range 64 - 133		Recovery = 92.820%			
Target Compounds						
16) Acetone	4.430	43	2794	3.799	ug/l	98
60) Dibromochloromethane	11.364	129	4330	1.880	ug/l	90
71) Bromoform	12.588	173	1913	1.079	ug/l #	99

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

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