

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN081524\
 Data File : VN083341.D
 Acq On : 15 Aug 2024 19:15
 Operator : JC\MD
 Sample : P3623-03
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 BPOW6-7-HYD-20240814

Quant Time: Aug 28 07:13:11 2024
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N080724W.M
 Quant Title : SW846 8260
 QLast Update : Thu Aug 08 06:30:41 2024
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.224	168	118987	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.100	114	237229	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.865	117	248301	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	108377	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.583	65	100457	59.314	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery = 118.620%			
35) Dibromofluoromethane	8.171	113	80260	54.203	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery = 108.400%			
50) Toluene-d8	10.571	98	292268	52.914	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery = 105.820%			
62) 4-Bromofluorobenzene	12.853	95	120978	56.181	ug/l	0.00
Spiked Amount 50.000	Range 77 - 121		Recovery = 112.360%			
Target Compounds						
30) Chloroform	7.971	83	1643	0.619	ug/l	94
47) Bromodichloromethane	9.882	83	3285	1.290	ug/l #	73
60) Dibromochloromethane	11.365	129	3685	2.017	ug/l	99
71) Bromoform	12.582	173	1205	0.822	ug/l #	89

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

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