

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN060923\
 Data File : VN078141.D
 Acq On : 09 Jun 2023 13:23
 Operator : JC\MD
 Sample : 03123-12
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 BP-VPB-RW11-EB-20230607

Quant Time: Jun 09 23:07:10 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

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

Internal Standards						
1) Pentafluorobenzene	8.230	168	489486	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	900231	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	800851	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.800	152	254790	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.583	65	349839	53.604	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	107.200%
35) Dibromofluoromethane	8.177	113	302230	53.097	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	106.200%
50) Toluene-d8	10.577	98	1082547	49.334	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	98.660%
62) 4-Bromofluorobenzene	12.853	95	347281	50.909	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	101.820%
Target Compounds						
						Qvalue
18) Methyl Acetate	5.048	43	11973	1.384	ug/l	90
30) Chloroform	7.971	83	10455	0.959	ug/l #	82

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

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