

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN091222\
 Data File : VN074418.D
 Acq On : 12 Sep 2022 19:11
 Operator : JC\MD
 Sample : N4558-01
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 BP-VPB-RW5-TB-20220906

Quant Time: Sep 13 06:27:51 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N090922W.M
 Quant Title : SW846 8260
 QLast Update : Sat Sep 10 06:10:29 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.016	168	296969	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.904	114	513723	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.686	117	511331	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.615	152	219726	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.375	65	226725	44.127	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery =	88.260%		
35) Dibromofluoromethane	7.957	113	188110	54.240	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery =	108.480%		
50) Toluene-d8	10.380	98	622052	46.032	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery =	92.060%		
62) 4-Bromofluorobenzene	12.674	95	243828	49.697	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery =	99.400%		
Target Compounds					Qvalue	
16) Acetone	4.228	43	37344	13.463	ug/l #	88
30) Chloroform	7.745	83	33770	4.027	ug/l #	79
47) Bromodichloromethane	9.698	83	4047	0.639	ug/l #	53

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

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