

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN081221\
 Data File : VN068103.D
 Acq On : 12 Aug 2021 20:46
 Operator : JC/MD
 Sample : M3361-09
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 41911

Quant Time: Aug 13 05:47:44 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N072921W.M
 Quant Title : SW846 8260
 QLast Update : Thu Jul 29 15:02:35 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.088	168	287941	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.971	114	561977	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.747	117	548718	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.680	152	186506	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.442	65	226409	55.543	ug/l	0.00
Spiked Amount 50.000	Range 61 - 141		Recovery = 111.080%			
35) Dibromofluoromethane	8.024	113	173473	52.967	ug/l	0.00
Spiked Amount 50.000	Range 69 - 133		Recovery = 105.940%			
50) Toluene-d8	10.446	98	713188	47.402	ug/l	0.00
Spiked Amount 50.000	Range 65 - 126		Recovery = 94.800%			
62) 4-Bromofluorobenzene	12.734	95	232917	42.886	ug/l	0.00
Spiked Amount 50.000	Range 58 - 135		Recovery = 85.780%			
Target Compounds						Qvalue
16) Acetone	4.293	43	4125	2.773	ug/l	93
30) Chloroform	7.820	83	64332	9.758	ug/l	99
47) Bromodichloromethane	9.765	83	26353	4.988	ug/l	98
60) Dibromochloromethane	11.240	129	12715	5.755	ug/l	94

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

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