

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN101921\
 Data File : VN069025.D
 Acq On : 19 Oct 2021 14:53
 Operator : JC/MD
 Sample : M4251-03RE
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 SW-RR-03-(SA)RE

Quant Time: Oct 20 05:48:31 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N101221W.M
 Quant Title : SW846 8260
 QLast Update : Wed Oct 13 05:44:40 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.088	168	334773	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.971	114	584110	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.746	117	530952	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.677	152	194526	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.440	65	215232	48.763	ug/l	0.00
Spiked Amount 50.000	Range 61 - 141		Recovery =	97.520%		
35) Dibromofluoromethane	8.024	113	171555	50.481	ug/l	0.00
Spiked Amount 50.000	Range 69 - 133		Recovery =	100.960%		
50) Toluene-d8	10.443	98	719469	52.912	ug/l	0.00
Spiked Amount 50.000	Range 65 - 126		Recovery =	105.820%		
62) 4-Bromofluorobenzene	12.733	95	245249	47.805	ug/l	0.00
Spiked Amount 50.000	Range 58 - 135		Recovery =	95.600%		
Target Compounds					Qvalue	
16) Acetone	4.304	43	4471	2.714	ug/l	99
30) Chloroform	7.825	83	2013	0.292	ug/l	89
52) Toluene	10.507	92	2915	0.298	ug/l	88

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

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