

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN101921\
 Data File : VN069040.D
 Acq On : 19 Oct 2021 21:10
 Operator : JC/MD
 Sample : M4259-01
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 WC-1-CS

Quant Time: Oct 20 05:50:54 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	340179	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.971	114	592056	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.747	117	565671	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.678	152	209209	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.440	65	222346	49.574	ug/l	0.00
Spiked Amount 50.000	Range 61 - 141		Recovery =	99.140%		
35) Dibromofluoromethane	8.024	113	174083	50.537	ug/l	0.00
Spiked Amount 50.000	Range 69 - 133		Recovery =	101.080%		
50) Toluene-d8	10.443	98	739446	53.651	ug/l	0.00
Spiked Amount 50.000	Range 65 - 126		Recovery =	107.300%		
62) 4-Bromofluorobenzene	12.734	95	259520	49.908	ug/l	0.00
Spiked Amount 50.000	Range 58 - 135		Recovery =	99.820%		
Target Compounds					Qvalue	
16) Acetone	4.299	43	16521	9.871	ug/l	99
18) Methyl Acetate	4.862	43	8996	1.417	ug/l #	70
20) Methylene Chloride	5.101	84	17520	4.548	ug/l	99
43) Isopropyl Acetate	8.566	43	55787	6.054	ug/l #	81

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

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