

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN121021\
 Data File : VN069987.D
 Acq On : 11 Dec 2021 5:08
 Operator : JC/MD
 Sample : M4990-20
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 R1-4

Quant Time: Dec 11 08:31:14 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N120221W.M
 Quant Title : SW846 8260
 QLast Update : Fri Dec 03 08:46:47 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.088	168	1063111	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.968	114	2092025	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.744	117	2440475	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.678	152	1032724	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.440	65	880475	53.640	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	= 107.280%	
35) Dibromofluoromethane	8.024	113	647786	49.434	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	= 98.860%	
50) Toluene-d8	10.443	98	2800434	53.223	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	= 106.440%	
62) 4-Bromofluorobenzene	12.731	95	1191620	62.513	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	= 125.020%	
Target Compounds						
					Qvalue	
16) Acetone	4.299	43	71765	7.268	ug/l	99
18) Methyl Acetate	4.870	43	46595	1.563	ug/l #	76
20) Methylene Chloride	5.101	84	56563	4.309	ug/l #	86
43) Isopropyl Acetate	8.566	43	151801	3.461	ug/l #	81

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN121021\
Data File : VN069987.D
Acq On : 11 Dec 2021 5:08
Operator : JC/MD
Sample : M4990-20
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 37 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
R1-4

Quant Time: Dec 11 08:31:14 2021
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N120221W.M
Quant Title : SW846 8260
QLast Update : Fri Dec 03 08:46:47 2021
Response via : Initial Calibration

