

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN120820\
 Data File : VN065003.D
 Acq On : 9 Dec 2020 5:51
 Operator : JC/MD
 Sample : L4970-16
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 49 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 4S1-A-BOTTOM

Quant Time: Dec 09 08:32:38 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N112320W.M
 Quant Title : SW846 8260
 QLast Update : Mon Nov 23 13:54:12 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.63	168	162279	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	314720	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	330043	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	137199	50.00	ug/l	0.00

System Monitoring Compounds

33) 1,2-Dichloroethane-d4	7.99	65	106045	56.21	ug/l	0.00
Spiked Amount 50.000	Range 61 - 141		Recovery =	112.42%		
35) Dibromofluoromethane	7.55	113	98585	52.35	ug/l	0.00
Spiked Amount 50.000	Range 69 - 133		Recovery =	104.70%		
50) Toluene-d8	10.06	98	393579	51.15	ug/l	0.00
Spiked Amount 50.000	Range 65 - 126		Recovery =	102.30%		
62) 4-Bromofluorobenzene	12.38	95	145261	52.75	ug/l	0.00
Spiked Amount 50.000	Range 58 - 135		Recovery =	105.50%		

Target Compounds

					Qvalue
16) Acetone	3.78	43	11636	18.762	ug/l 99
20) Methylene Chloride	4.50	84	18611	9.577	ug/l 89
43) Isopropyl Acetate	8.13	43	15236	3.370	ug/l # 93

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN120820\
Data File : VN065003.D
Acq On : 9 Dec 2020 5:51
Operator : JC/MD
Sample : L4970-16
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 49 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
4S1-A-BOTTOM

Quant Time: Dec 09 08:32:38 2020
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N112320W.M
Quant Title : SW846 8260
QLast Update : Mon Nov 23 13:54:12 2020
Response via : Initial Calibration

