

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN031622\
 Data File : VN071328.D
 Acq On : 16 Mar 2022 18:25
 Operator : JC\MD
 Sample : N1957-10
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 MP-1D

Quant Time: Mar 17 05:20:20 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N030722W.M
 Quant Title : SW846 8260
 QLast Update : Tue Mar 08 06:52:46 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.086	168	903292	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.963	114	1476446	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.745	117	1375539	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.668	152	523447	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.434	65	560841	45.302	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	=	90.600%
35) Dibromofluoromethane	8.022	113	429908	46.143	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	=	92.280%
50) Toluene-d8	10.439	98	1805482	48.130	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	=	96.260%
62) 4-Bromofluorobenzene	12.727	95	675196	53.154	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	=	106.300%
Target Compounds						
					Qvalue	
16) Acetone	4.293	43	14598	2.423	ug/l	93
18) Methyl Acetate	4.869	43	17648	1.551	ug/l	92
64) Tetrachloroethene	10.975	164	30759	3.206	ug/l	98

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN031622\
Data File : VN071328.D
Acq On : 16 Mar 2022 18:25
Operator : JC\MD
Sample : N1957-10
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MP-1D

Quant Time: Mar 17 05:20:20 2022
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N030722W.M
Quant Title : SW846 8260
QLast Update : Tue Mar 08 06:52:46 2022
Response via : Initial Calibration

