

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN120721\
 Data File : VN069813.D
 Acq On : 7 Dec 2021 14:33
 Operator : JC/MD
 Sample : M4956-02
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 RE134D1-20211202

Quant Time: Dec 08 04:35:33 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	8.086	168	1302794	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.968	114	2343837	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.744	117	2454151	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.675	152	949559	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.440	65	1004432	49.934	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	=	99.860%
35) Dibromofluoromethane	8.024	113	694782	47.324	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	=	94.640%
50) Toluene-d8	10.443	98	2978440	50.525	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	=	101.040%
62) 4-Bromofluorobenzene	12.731	95	1159353	54.286	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	=	108.580%
Target Compounds						
					Qvalue	
16) Acetone	4.296	43	21884	1.809	ug/l	98
44) Trichloroethene	9.217	130	45580	2.800	ug/l	94
64) Tetrachloroethene	10.977	164	6530	0.392	ug/l	88

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN120721\
Data File : VN069813.D
Acq On : 7 Dec 2021 14:33
Operator : JC/MD
Sample : M4956-02
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
RE134D1-20211202

Quant Time: Dec 08 04:35:33 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

