

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN070722\
 Data File : VN073374.D
 Acq On : 07 Jul 2022 16:21
 Operator : JC\MD
 Sample : N3620-02
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 BP-TT-MW174I1-GW-20220706

Quant Time: Jul 08 03:02:13 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N062822W.M
 Quant Title : SW846 8260
 QLast Update : Wed Jun 29 01:54:03 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.016	168	290529	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.904	114	486486	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.686	117	451105	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.610	152	209601	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.369	65	182393	43.488	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	86.980%
35) Dibromofluoromethane	7.951	113	154469	47.725	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	95.440%
50) Toluene-d8	10.381	98	516873	44.963	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	89.920%
62) 4-Bromofluorobenzene	12.669	95	223578	50.610	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	101.220%
Target Compounds						
					Qvalue	
3) Chloromethane	2.269	50	1301	0.315	ug/l	89
19) Methyl tert-butyl Ether	5.552	73	4628	0.371	ug/l	91
24) 1,1-Dichloroethane	6.304	63	25463	3.660	ug/l #	95

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN070722\
Data File : VN073374.D
Acq On : 07 Jul 2022 16:21
Operator : JC\MD
Sample : N3620-02
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
BP-TT-MW174I1-GW-20220706

Quant Time: Jul 08 03:02:13 2022
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N062822W.M
Quant Title : SW846 8260
QLast Update : Wed Jun 29 01:54:03 2022
Response via : Initial Calibration

