

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN090922\
 Data File : VN074387.D
 Acq On : 09 Sep 2022 20:58
 Operator : JC\MD
 Sample : N4471-10
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 MW-14A-15-20-083122

Quant Time: Sep 10 07:26:47 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N090922W.M
 Quant Title : SW846 8260
 QLast Update : Sat Sep 10 06:10:29 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.016	168	341290	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.904	114	593622	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.686	117	586161	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.616	152	277329	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.369	65	242417	41.054	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery =	82.100%		
35) Dibromofluoromethane	7.951	113	202012	50.408	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery =	100.820%		
50) Toluene-d8	10.381	98	678434	43.447	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery =	86.900%		
62) 4-Bromofluorobenzene	12.674	95	279877	49.367	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery =	98.740%		
Target Compounds						Qvalue
27) cis-1,2-Dichloroethene	7.251	96	3925	0.668	ug/l	81
30) Chloroform	7.751	83	6764	0.702	ug/l	92
44) Trichloroethene	9.145	130	12808	2.826	ug/l	97

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN090922\
Data File : VN074387.D
Acq On : 09 Sep 2022 20:58
Operator : JC\MD
Sample : N4471-10
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-14A-15-20-083122

Quant Time: Sep 10 07:26:47 2022
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N090922W.M
Quant Title : SW846 8260
QLast Update : Sat Sep 10 06:10:29 2022
Response via : Initial Calibration

