

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN110423\
 Data File : VN079897.D
 Acq On : 04 Nov 2023 16:53
 Operator : JC\MD
 Sample : 05222-07DL 5X
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 FSND-MW-EVAL-02S-20231101DL

Quant Time: Nov 06 02:16:14 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N110123W.M
 Quant Title : SW846 8260
 QLast Update : Thu Nov 02 05:49:08 2023
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.230	168	228619	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	464160	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	475496	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	181834	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.583	65	205311	58.945	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	117.900%
35) Dibromofluoromethane	8.171	113	169009	51.447	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	102.900%
50) Toluene-d8	10.571	98	579526	47.919	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	95.840%
62) 4-Bromofluorobenzene	12.853	95	190976	47.614	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	95.220%
Target Compounds						
						Qvalue
44) Trichloroethene	9.359	130	80408	23.065	ug/l	96
64) Tetrachloroethene	11.112	164	16626	5.399	ug/l	93

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN110423\
Data File : VN079897.D
Acq On : 04 Nov 2023 16:53
Operator : JC\MD
Sample : 05222-07DL 5X
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
FSND-MW-EVAL-02S-20231101DL

Quant Time: Nov 06 02:16:14 2023
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N110123W.M
Quant Title : SW846 8260
QLast Update : Thu Nov 02 05:49:08 2023
Response via : Initial Calibration

