

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN111124\
 Data File : VN084779.D
 Acq On : 11 Nov 2024 18:06
 Operator : JC\MD
 Sample : P4788-04
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 BP-G3

Quant Time: Nov 12 00:36:10 2024
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N103024W.M
 Quant Title : SW846 8260
 QLast Update : Thu Oct 31 18:45:38 2024
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.224	168	164711	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.100	114	297853	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.865	117	276056	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	122850	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.577	65	117202	49.266	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	98.540%
35) Dibromofluoromethane	8.165	113	95717	47.476	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	94.960%
50) Toluene-d8	10.565	98	354051	47.673	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	95.340%
62) 4-Bromofluorobenzene	12.847	95	135026	48.648	ug/l	0.00
Spiked Amount	50.000	Range	77 - 121	Recovery	=	97.300%
Target Compounds						
16) Acetone	4.430	43	23962	32.386	ug/l	99
20) Methylene Chloride	5.271	84	3525	1.685	ug/l #	63
43) Isopropyl Acetate	8.688	43	206580	43.962	ug/l #	73

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN111124\
Data File : VN084779.D
Acq On : 11 Nov 2024 18:06
Operator : JC\MD
Sample : P4788-04
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
BP-G3

Quant Time: Nov 12 00:36:10 2024
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N103024W.M
Quant Title : SW846 8260
QLast Update : Thu Oct 31 18:45:38 2024
Response via : Initial Calibration

