

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN111824\
 Data File : VN084930.D
 Acq On : 18 Nov 2024 19:16
 Operator : JC\MD
 Sample : P4870-15
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 MH-736

Quant Time: Nov 19 00:34:21 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	8.218	168	198291	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.100	114	363891	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.865	117	327811	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.788	152	151104	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.571	65	145704	50.874	ug/l	0.00
Spiked Amount	50.000	Range 74 - 125	Recovery	= 101.740%		
35) Dibromofluoromethane	8.159	113	113231	45.971	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	= 91.940%		
50) Toluene-d8	10.565	98	429627	47.351	ug/l	0.00
Spiked Amount	50.000	Range 86 - 113	Recovery	= 94.700%		
62) 4-Bromofluorobenzene	12.847	95	174769	51.539	ug/l	0.00
Spiked Amount	50.000	Range 77 - 121	Recovery	= 103.080%		
Target Compounds						
					Qvalue	
16) Acetone	4.442	43	19038	20.392	ug/l	95
43) Isopropyl Acetate	8.688	43	233416	40.584	ug/l #	76
64) Tetrachloroethene	11.100	164	2779	1.274	ug/l #	75

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN111824\
Data File : VN084930.D
Acq On : 18 Nov 2024 19:16
Operator : JC\MD
Sample : P4870-15
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MH-736

Quant Time: Nov 19 00:34:21 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

