

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN111423\
 Data File : VN080094.D
 Acq On : 14 Nov 2023 17:14
 Operator : JC\MD
 Sample : 05378-04
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 MH-3

Quant Time: Nov 15 03:57:39 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	182173	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	358577	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	386282	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	163505	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.583	65	162779	58.649	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	117.300%
35) Dibromofluoromethane	8.165	113	126284	49.760	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	99.520%
50) Toluene-d8	10.571	98	466899	49.974	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	99.940%
62) 4-Bromofluorobenzene	12.853	95	182503	58.900	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	117.800%
Target Compounds						
					Qvalue	
16) Acetone	4.430	43	22790	23.377	ug/l	99
20) Methylene Chloride	5.277	84	7969	3.239	ug/l	90
43) Isopropyl Acetate	8.688	43	129130	22.157	ug/l #	84

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN111423\
Data File : VN080094.D
Acq On : 14 Nov 2023 17:14
Operator : JC\MD
Sample : 05378-04
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MH-3

Quant Time: Nov 15 03:57:39 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

