

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN042222\
 Data File : VN072155.D
 Acq On : 22 Apr 2022 14:34
 Operator : JC\MD
 Sample : N2526-08
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 MH-5D

Quant Time: Apr 23 03:17:41 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N033022W.M
 Quant Title : SW846 8260
 QLast Update : Thu Mar 31 01:57:54 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.081	168	512558	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.963	114	872755	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.739	117	969374	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.669	152	355906	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.434	65	272526	46.574	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	=	93.140%
35) Dibromofluoromethane	8.016	113	249438	48.570	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	=	97.140%
50) Toluene-d8	10.439	98	1136965	54.118	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	=	108.240%
62) 4-Bromofluorobenzene	12.727	95	391364	57.599	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	=	115.200%
Target Compounds						
					Qvalue	
16) Acetone	4.293	43	24060	9.621	ug/l	94
20) Methylene Chloride	5.099	84	41385	7.852	ug/l	90
43) Isopropyl Acetate	8.557	43	43690	3.346	ug/l #	92

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN042222\
Data File : VN072155.D
Acq On : 22 Apr 2022 14:34
Operator : JC\MD
Sample : N2526-08
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MH-5D

Quant Time: Apr 23 03:17:41 2022
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N033022W.M
Quant Title : SW846 8260
QLast Update : Thu Mar 31 01:57:54 2022
Response via : Initial Calibration

