

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN042122\
 Data File : VN072131.D
 Acq On : 21 Apr 2022 17:33
 Operator : JC\MD
 Sample : N2481-01
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 MW-12

Quant Time: Apr 22 02:47:34 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	8.086	168	551504	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.963	114	943591	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.739	117	1001690	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.668	152	391488	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.433	65	299592	47.584	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	=	95.160%
35) Dibromofluoromethane	8.022	113	267184	48.120	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	=	96.240%
50) Toluene-d8	10.439	98	1189929	52.388	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	=	104.780%
62) 4-Bromofluorobenzene	12.727	95	416208	56.657	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	=	113.320%
Target Compounds						
					Qvalue	
16) Acetone	4.293	43	27735	10.308	ug/l #	88
25) 2-Butanone	7.345	43	3179	0.820	ug/l #	84
37) Ethyl Acetate	7.416	43	12585	1.483	ug/l	99

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN042122\
Data File : VN072131.D
Acq On : 21 Apr 2022 17:33
Operator : JC\MD
Sample : N2481-01
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-12

Quant Time: Apr 22 02:47:34 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

