

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN030322\
 Data File : VN071157.D
 Acq On : 03 Mar 2022 16:46
 Operator : JC\MD
 Sample : N1750-07
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 WC-2-Q4

Quant Time: Mar 04 01:15:20 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N021522W.M
 Quant Title : SW846 8260
 QLast Update : Mon Feb 21 07:57:15 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.081	168	872443	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.963	114	1426732	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.739	117	1287850	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.675	152	446891	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.434	65	531257	53.307	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	=	106.620%
35) Dibromofluoromethane	8.022	113	410043	48.319	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	=	96.640%
50) Toluene-d8	10.440	98	1747310	51.271	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	=	102.540%
62) 4-Bromofluorobenzene	12.728	95	593175	48.694	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	=	97.380%
Target Compounds						
						Qvalue
16) Acetone	4.287	43	47090	8.580	ug/l	91
18) Methyl Acetate	4.858	43	26420	1.615	ug/l	91
20) Methylene Chloride	5.093	84	19487	1.954	ug/l	97
43) Isopropyl Acetate	8.557	43	132341	5.631	ug/l #	77

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN030322\
Data File : VN071157.D
Acq On : 03 Mar 2022 16:46
Operator : JC\MD
Sample : N1750-07
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
WC-2-Q4

Quant Time: Mar 04 01:15:20 2022
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N021522W.M
Quant Title : SW846 8260
QLast Update : Mon Feb 21 07:57:15 2022
Response via : Initial Calibration

