

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN030322\
 Data File : VN071155.D
 Acq On : 03 Mar 2022 15:58
 Operator : JC\MD
 Sample : N1750-05
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 WC-2-Q2

Quant Time: Mar 04 01:14:44 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	752358	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.963	114	1241922	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.739	117	1120163	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.674	152	384660	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.434	65	472658	54.997	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	=	110.000%
35) Dibromofluoromethane	8.022	113	358931	48.590	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	=	97.180%
50) Toluene-d8	10.439	98	1531050	51.611	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	=	103.220%
62) 4-Bromofluorobenzene	12.727	95	511592	48.247	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	=	96.500%
Target Compounds						
						Qvalue
16) Acetone	4.293	43	44594	9.422	ug/l	99
18) Methyl Acetate	4.863	43	23757	1.710	ug/l	94
20) Methylene Chloride	5.104	84	19185	2.230	ug/l #	87
43) Isopropyl Acetate	8.557	43	134529	6.576	ug/l #	77

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN030322\
Data File : VN071155.D
Acq On : 03 Mar 2022 15:58
Operator : JC\MD
Sample : N1750-05
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
WC-2-Q2

Quant Time: Mar 04 01:14:44 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

