

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN012422\
 Data File : VN070759.D
 Acq On : 24 Jan 2022 16:45
 Operator : JC/MD
 Sample : N1235-02
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 TP-1

Quant Time: Jan 24 23:33:20 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N011822W.M
 Quant Title : SW846 8260
 QLast Update : Tue Jan 18 13:32:07 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.086	168	946475	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.965	114	1547513	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.744	117	1382967	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.675	152	503596	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.437	65	644525	47.516	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	=	95.040%
35) Dibromofluoromethane	8.021	113	443732	46.226	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	=	92.460%
50) Toluene-d8	10.440	98	1936647	47.890	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	=	95.780%
62) 4-Bromofluorobenzene	12.731	95	624861	43.362	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	=	86.720%
Target Compounds						
					Qvalue	
16) Acetone	4.288	43	99205	10.537	ug/l	94
18) Methyl Acetate	4.862	43	27366	1.478	ug/l #	76
20) Methylene Chloride	5.093	84	26553	2.141	ug/l #	88
43) Isopropyl Acetate	8.563	43	137796	4.205	ug/l #	83

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN012422\
Data File : VN070759.D
Acq On : 24 Jan 2022 16:45
Operator : JC/MD
Sample : N1235-02
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
TP-1

Quant Time: Jan 24 23:33:20 2022
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N011822W.M
Quant Title : SW846 8260
QLast Update : Tue Jan 18 13:32:07 2022
Response via : Initial Calibration

