

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN031522\
 Data File : VN071299.D
 Acq On : 15 Mar 2022 18:04
 Operator : JC\MD
 Sample : N1918-02
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 AVENUE-X-GROUT-SAMPLE-A

Quant Time: Mar 16 08:05:17 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N030722W.M
 Quant Title : SW846 8260
 QLast Update : Tue Mar 08 06:52:46 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.086	168	811191	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.963	114	1370552	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.739	117	1253924	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.674	152	449740	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.433	65	519729	46.748	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	=	93.500%
35) Dibromofluoromethane	8.022	113	247638	28.633	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	=	57.260%#
50) Toluene-d8	10.439	98	1651751	47.434	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	=	94.860%
62) 4-Bromofluorobenzene	12.727	95	601557	51.016	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	=	102.040%
Target Compounds						
					Qvalue	
16) Acetone	4.281	43	244322	45.165	ug/l	97
18) Methyl Acetate	4.863	43	19975	1.832	ug/l	95
20) Methylene Chloride	5.098	84	15390	1.927	ug/l	95
25) 2-Butanone	7.334	43	37307	4.281	ug/l	100

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN031522\
Data File : VN071299.D
Acq On : 15 Mar 2022 18:04
Operator : JC\MD
Sample : N1918-02
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
AVENUE-X-GROUT-SAMPLE-A

Quant Time: Mar 16 08:05:17 2022
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N030722W.M
Quant Title : SW846 8260
QLast Update : Tue Mar 08 06:52:46 2022
Response via : Initial Calibration

