

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN031422\
 Data File : VN071268.D
 Acq On : 14 Mar 2022 16:02
 Operator : JC\MD
 Sample : N1918-05
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 13 Sample Multiplier: 1

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

Quant Time: Mar 15 02:52:53 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	8.081	168	788020	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.963	114	1300953	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.739	117	1179593	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.668	152	437401	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.433	65	494435	45.781	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	=	91.560%
35) Dibromofluoromethane	8.016	113	234553	28.571	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	=	57.140%#
50) Toluene-d8	10.439	98	1573915	47.617	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	=	95.240%
62) 4-Bromofluorobenzene	12.727	95	573366	51.226	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	=	102.460%
Target Compounds						
					Qvalue	
16) Acetone	4.281	43	258684	49.226	ug/l	98
18) Methyl Acetate	4.851	43	25945	2.291	ug/l	95
20) Methylene Chloride	5.098	84	14733	1.904	ug/l	96
25) 2-Butanone	7.334	43	42514	5.022	ug/l	96

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN031422\
Data File : VN071268.D
Acq On : 14 Mar 2022 16:02
Operator : JC\MD
Sample : N1918-05
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 13 Sample Multiplier: 1

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

Quant Time: Mar 15 02:52:53 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

