

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN031422\
 Data File : VN071267.D
 Acq On : 14 Mar 2022 15:38
 Operator : JC\MD
 Sample : N1918-02
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 12 Sample Multiplier: 1

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

Quant Time: Mar 15 02:52:39 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	8.080	168	717538	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.963	114	1184476	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.739	117	1067608	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.668	152	389819	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.433	65	459759	46.751	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	=	93.500%
35) Dibromofluoromethane	8.016	113	213679	28.588	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	=	57.180%#
50) Toluene-d8	10.439	98	1427757	47.443	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	=	94.880%
62) 4-Bromofluorobenzene	12.727	95	502301	49.290	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	=	98.580%
Target Compounds						
						Qvalue
16) Acetone	4.281	43	254638	53.216	ug/l	97
18) Methyl Acetate	4.863	43	31237	2.878	ug/l	97
20) Methylene Chloride	5.092	84	15632	2.156	ug/l	95
25) 2-Butanone	7.333	43	43231	5.609	ug/l	93

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

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