

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN120321\
 Data File : VN069743.D
 Acq On : 4 Dec 2021 2:41
 Operator : JC/MD
 Sample : M4919-01
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 38 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 FDR-MW-2-20211202

Quant Time: Dec 04 05:41:05 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N120221W.M
 Quant Title : SW846 8260
 QLast Update : Fri Dec 03 08:46:47 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.086	168	1515986	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.968	114	2664542	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.744	117	2744331	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.675	152	1064223	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.437	65	1161811	49.636	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	=	99.280%
35) Dibromofluoromethane	8.021	113	799927	47.928	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	=	95.860%
50) Toluene-d8	10.440	98	3398013	50.704	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	=	101.400%
62) 4-Bromofluorobenzene	12.731	95	1273813	52.467	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	=	104.940%
Target Compounds						
					Qvalue	
16) Acetone	4.299	43	27966	1.986	ug/l	99
17) Carbon Disulfide	4.537	76	21517	0.577	ug/l	96
24) 1,1-Dichloroethane	6.385	63	9474	0.288	ug/l #	94
83) tert-Butylbenzene	13.323	119	386129	5.722	ug/l	92

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

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