

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN110221\
 Data File : VN069343.D
 Acq On : 2 Nov 2021 12:59
 Operator : JC/MD
 Sample : M4438-04
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 FDR-20211101-WC-LS

Quant Time: Nov 03 00:47:48 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N102821W.M
 Quant Title : SW846 8260
 QLast Update : Thu Oct 28 16:09:15 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.088	168	289052	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.971	114	518976	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.746	117	543273	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.677	152	223679	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.440	65	204674	49.052	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	=	98.100%
35) Dibromofluoromethane	8.024	113	152541	47.152	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	=	94.300%
50) Toluene-d8	10.443	98	659429	50.980	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	=	101.960%
62) 4-Bromofluorobenzene	12.733	95	261437	53.562	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	=	107.120%
Target Compounds						
					Qvalue	
16) Acetone	4.296	43	23921	13.456	ug/l	97
17) Carbon Disulfide	4.527	76	5503	0.795	ug/l	96
25) 2-Butanone	7.353	43	7102	4.224	ug/l #	87
40) Benzene	8.464	78	5921	0.440	ug/l	95
56) Ethyl methacrylate	10.762	69	6741	2.685	ug/l	97
83) tert-Butylbenzene	13.323	119	66446	4.879	ug/l	98

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

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