

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN110221\
 Data File : VN069341.D
 Acq On : 2 Nov 2021 12:08
 Operator : JC/MD
 Sample : M4438-02
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 7 Sample Multiplier: 1

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

Quant Time: Nov 03 00:47:23 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	276962	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.971	114	489420	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.746	117	494386	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.677	152	200479	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.440	65	189529	47.405	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	=	94.820%
35) Dibromofluoromethane	8.024	113	144330	47.308	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	=	94.620%
50) Toluene-d8	10.443	98	616119	50.509	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	=	101.020%
62) 4-Bromofluorobenzene	12.733	95	237759	51.652	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	=	103.300%
Target Compounds						
					Qvalue	
16) Acetone	4.299	43	33901	19.902	ug/l	97
17) Carbon Disulfide	4.529	76	12904	1.945	ug/l	99
25) 2-Butanone	7.345	43	9280	5.166	ug/l	95
40) Benzene	8.464	78	6625	0.522	ug/l	95
56) Ethyl methacrylate	10.762	69	5350	2.526	ug/l	94
83) tert-Butylbenzene	13.323	119	70944	5.812	ug/l	97

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

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