

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN072321\
 Data File : VN067887.D
 Acq On : 23 Jul 2021 18:06
 Operator : JC/MD
 Sample : M3147-08
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 152140-MW-2D(ONSITE)

Manual Integrations
 APPROVED

MMDadoda
 7/26/2021 5:22:01 PM

Quant Time: Jul 24 01:43:16 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N070721W.M
 Quant Title : SW846 8260
 QLast Update : Wed Jul 07 16:18:53 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.088	168	268087	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.971	114	514121	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.749	117	483344	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.678	152	161574	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.440	65	214543	52.952	ug/l	0.00
Spiked Amount 50.000	Range 61 - 141		Recovery = 105.900%			
35) Dibromofluoromethane	8.027	113	157406	50.675	ug/l	0.00
Spiked Amount 50.000	Range 69 - 133		Recovery = 101.340%			
50) Toluene-d8	10.443	98	646421	49.718	ug/l	0.00
Spiked Amount 50.000	Range 65 - 126		Recovery = 99.440%			
62) 4-Bromofluorobenzene	12.733	95	201407	43.851	ug/l	0.00
Spiked Amount 50.000	Range 58 - 135		Recovery = 87.700%			
Target Compounds						
16) Acetone	4.309	43	3802m	2.982	ug/l	
44) Trichloroethene	9.217	130	4481	1.235	ug/l	83
64) Tetrachloroethene	10.979	164	47482	13.917	ug/l	97

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

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