

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN072621\
 Data File : VN067910.D
 Acq On : 26 Jul 2021 16:27
 Operator : JC/MD
 Sample : M3147-13
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 152140-MW-4S(ONSITE)

Manual Integrations
 APPROVED

MMDadoda
 7/27/2021 3:56:17 PM

Quant Time: Jul 27 01:30:52 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	8.091	168	309669	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.971	114	585548	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.749	117	557359	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.677	152	181717	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.440	65	243998	52.136	ug/l	0.00
Spiked Amount 50.000	Range 61 - 141		Recovery = 104.280%			
35) Dibromofluoromethane	8.024	113	179504	50.740	ug/l	0.00
Spiked Amount 50.000	Range 69 - 133		Recovery = 101.480%			
50) Toluene-d8	10.446	98	754578	50.957	ug/l	0.00
Spiked Amount 50.000	Range 65 - 126		Recovery = 101.920%			
62) 4-Bromofluorobenzene	12.733	95	235072	44.938	ug/l	0.00
Spiked Amount 50.000	Range 58 - 135		Recovery = 89.880%			
Target Compounds					Qvalue	
16) Acetone	4.304	43	5509	3.741	ug/l	98
18) Methyl Acetate	4.867	43	4138m	1.000	ug/l	
44) Trichloroethene	9.215	130	2058	0.498	ug/l	91
64) Tetrachloroethene	10.977	164	10809	2.747	ug/l	88

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

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