

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN102121\
 Data File : VN069090.D
 Acq On : 21 Oct 2021 16:02
 Operator : JC/MD
 Sample : M4284-13 200X
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 30856

Manual Integrations
 APPROVED

MMDadoda
 10/22/2021 7:09:18 PM

Quant Time: Oct 22 01:33:43 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N101221W.M
 Quant Title : SW846 8260
 QLast Update : Wed Oct 13 05:44:40 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.088	168	315326	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.971	114	555452	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.746	117	514232	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.677	152	207260	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.442	65	209678	50.434	ug/l	0.00
Spiked Amount 50.000	Range 61 - 141		Recovery = 100.860%			
35) Dibromofluoromethane	8.024	113	163587	50.620	ug/l	0.00
Spiked Amount 50.000	Range 69 - 133		Recovery = 101.240%			
50) Toluene-d8	10.446	98	691246	53.459	ug/l	0.00
Spiked Amount 50.000	Range 65 - 126		Recovery = 106.920%			
62) 4-Bromofluorobenzene	12.733	95	246235	50.473	ug/l	0.00
Spiked Amount 50.000	Range 58 - 135		Recovery = 100.940%			
Target Compounds					Qvalue	
16) Acetone	4.299	43	6675	4.302	ug/l #	88
20) Methylene Chloride	5.103	84	3647m	1.021	ug/l	
30) Chloroform	7.823	83	154495	23.754	ug/l	99
37) Ethyl Acetate	7.423	43	31299m	6.388	ug/l	
86) p-Isopropyltoluene	13.618	119	28434	2.046	ug/l	99

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

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