

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN102121\
 Data File : VN069089.D
 Acq On : 21 Oct 2021 15:37
 Operator : JC/MD
 Sample : M4284-12 200X
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 30855

Manual Integrations
 APPROVED

MMDadoda
 10/22/2021 7:09:14 PM

Quant Time: Oct 22 01:33:33 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	329570	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.971	114	584254	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.749	117	550829	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.678	152	211471	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.440	65	220246	50.686	ug/l	0.00
Spiked Amount 50.000	Range 61 - 141		Recovery = 101.380%			
35) Dibromofluoromethane	8.027	113	172303	50.689	ug/l	0.00
Spiked Amount 50.000	Range 69 - 133		Recovery = 101.380%			
50) Toluene-d8	10.443	98	727911	53.519	ug/l	0.00
Spiked Amount 50.000	Range 65 - 126		Recovery = 107.040%			
62) 4-Bromofluorobenzene	12.733	95	260974	50.858	ug/l	0.00
Spiked Amount 50.000	Range 58 - 135		Recovery = 101.720%			
Target Compounds						
16) Acetone	4.312	43	6014m	3.709	ug/l	
20) Methylene Chloride	5.103	84	4961	1.329	ug/l	85
30) Chloroform	7.817	83	28241	4.155	ug/l	97
86) p-Isopropyltoluene	13.618	119	31386	2.214	ug/l	92

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

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