

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN120821\
 Data File : VN069897.D
 Acq On : 9 Dec 2021 7:18
 Operator : JC/MD
 Sample : M4959-20
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 49 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 TP-3D

Manual Integrations
 APPROVED

Reviewed By : John Carlone 12/10/2021
 Supervised By : Mahesh Dadoda 12/15/2021

Quant Time: Dec 09 12:01:38 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N120221W.M
 Quant Title : SW846 8260
 QLast Update : Fri Dec 03 08:46:47 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.088	168	1331473	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.968	114	2483538	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.744	117	2753012	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.675	152	1132193	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.440	65	1067727	51.937	ug/l	0.00
Spiked Amount 50.000	Range 61 - 141		Recovery = 103.880%			
35) Dibromofluoromethane	8.024	113	757623	48.702	ug/l	0.00
Spiked Amount 50.000	Range 69 - 133		Recovery = 97.400%			
50) Toluene-d8	10.440	98	3279454	52.502	ug/l	0.00
Spiked Amount 50.000	Range 65 - 126		Recovery = 105.000%			
62) 4-Bromofluorobenzene	12.731	95	1331947	58.859	ug/l	0.00
Spiked Amount 50.000	Range 58 - 135		Recovery = 117.720%			
Target Compounds						
16) Acetone	4.291	43	93470	7.558	ug/l	94
18) Methyl Acetate	4.870	43	64808m	1.736	ug/l	
20) Methylene Chloride	5.098	84	31412	1.911	ug/l	87
43) Isopropyl Acetate	8.563	43	196944	3.782	ug/l #	79

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

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