

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN051921\
 Data File : VN067154.D
 Acq On : 19 May 2021 16:16
 Operator : JC/MD
 Sample : M2366-09
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 JG-SPOILS-5-11-21-CRE

Manual Integrations
 APPROVED

SAM
 5/20/2021 6:04:19 PM

Quant Time: May 20 01:56:55 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N051421W.M
 Quant Title : SW846 8260
 QLast Update : Tue May 18 04:00:15 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	7.597	168	633641	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.525	114	873314	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.360	117	745393	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.307	152	269799	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.962	65	303651	44.36	ug/l	0.00
Spiked Amount 50.000	Range 61 - 141		Recovery =	88.72%		
35) Dibromofluoromethane	7.522	113	179452	31.29	ug/l	0.00
Spiked Amount 50.000	Range 69 - 133		Recovery =	62.58%#		
50) Toluene-d8	10.038	98	1084910	49.92	ug/l	0.00
Spiked Amount 50.000	Range 65 - 126		Recovery =	99.84%		
62) 4-Bromofluorobenzene	12.360	95	306218	42.06	ug/l	0.00
Spiked Amount 50.000	Range 58 - 135		Recovery =	84.12%		
Target Compounds					Qvalue	
16) Acetone	3.773	43	128521	60.41	ug/l	98
20) Methylene Chloride	4.489	84	34472	5.02	ug/l	94
25) 2-Butanone	6.768	43	44162	13.31	ug/l	98
51) 4-Methyl-2-Pentanone	9.938	43	8218m	6.55	ug/l	
69) o-Xylene	11.907	106	16784	1.54	ug/l	92

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

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