

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN071921\
 Data File : VN067806.D
 Acq On : 19 Jul 2021 20:28
 Operator : JC/MD
 Sample : M3089-04
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PAR-N2-A-TOP-(0)

Quant Time: Jul 20 04:55:02 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	8.088	168	310725	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.971	114	596722	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.749	117	611353	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.678	152	221797	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.440	65	247894	52.788	ug/l	0.00
Spiked Amount 50.000	Range 61 - 141		Recovery = 105.580%			
35) Dibromofluoromethane	8.024	113	180297	50.010	ug/l	0.00
Spiked Amount 50.000	Range 69 - 133		Recovery = 100.020%			
50) Toluene-d8	10.443	98	781147	51.764	ug/l	0.00
Spiked Amount 50.000	Range 65 - 126		Recovery = 103.520%			
62) 4-Bromofluorobenzene	12.734	95	265931	49.885	ug/l	0.00
Spiked Amount 50.000	Range 58 - 135		Recovery = 99.780%			
Target Compounds					Qvalue	
16) Acetone	4.291	43	35344	23.919	ug/l	93
18) Methyl Acetate	4.867	43	6220	1.499	ug/l #	77
20) Methylene Chloride	5.093	84	32339	8.802	ug/l	87
43) Isopropyl Acetate	8.566	43	36260	3.838	ug/l #	93

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

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