

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN051821\
 Data File : VN067138.D
 Acq On : 18 May 2021 19:39
 Operator : JC/MD
 Sample : M2366-07
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 22 Sample Multiplier: 1

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

Quant Time: May 19 02:52:49 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	617667	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.525	114	851216	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.357	117	725203	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.302	152	299993	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.959	65	277996	41.66	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	=	83.32%
35) Dibromofluoromethane	7.519	113	185999	33.28	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	=	66.56%#
50) Toluene-d8	10.038	98	993581	46.90	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	=	93.80%
62) 4-Bromofluorobenzene	12.358	95	302109	42.57	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	=	85.14%
Target Compounds						
					Qvalue	
16) Acetone	3.770	43	125242	60.39	ug/l	99
18) Methyl Acetate	4.274	43	23875	3.87	ug/l	98
20) Methylene Chloride	4.486	84	32353	4.84	ug/l	99
25) 2-Butanone	6.760	43	38681	11.96	ug/l	96
51) 4-Methyl-2-Pentanone	9.936	43	5864	6.24	ug/l	96

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

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