

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN051821\
 Data File : VN067137.D
 Acq On : 18 May 2021 19:15
 Operator : JC/MD
 Sample : M2366-09
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 21 Sample Multiplier: 1

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

Quant Time: May 19 02:52:33 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	613723	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.525	114	846210	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.360	117	723022	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.304	152	273955	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.962	65	276985	41.78	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	=	83.56%
35) Dibromofluoromethane	7.519	113	185242	33.34	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	=	66.68%#
50) Toluene-d8	10.037	98	987115	46.87	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	=	93.74%
62) 4-Bromofluorobenzene	12.360	95	293790	41.65	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	=	83.30%
Target Compounds						
					Qvalue	
16) Acetone	3.772	43	118830	57.67	ug/l	98
18) Methyl Acetate	4.277	43	23969	3.91	ug/l	97
20) Methylene Chloride	4.488	84	33458	5.03	ug/l	93
25) 2-Butanone	6.763	43	42499	13.23	ug/l	95
51) 4-Methyl-2-Pentanone	9.941	43	6540	6.34	ug/l	99
69) o-Xylene	11.909	106	18875	1.79	ug/l	92

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

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