

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN051821\
 Data File : VN067124.D
 Acq On : 18 May 2021 13:57
 Operator : JC/MD
 Sample : M2387-09 10X
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 30794

Quant Time: May 19 02:48:44 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	7.592	168	620840	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.522	114	870233	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.357	117	809536	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.302	152	326234	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.956	65	289223	43.12	ug/l	0.00
Spiked Amount 50.000	Range 61 - 141		Recovery =	86.24%		
35) Dibromofluoromethane	7.514	113	261945	45.84	ug/l	0.00
Spiked Amount 50.000	Range 69 - 133		Recovery =	91.68%		
50) Toluene-d8	10.035	98	1054044	48.67	ug/l	0.00
Spiked Amount 50.000	Range 65 - 126		Recovery =	97.34%		
62) 4-Bromofluorobenzene	12.358	95	350089	48.26	ug/l	0.00
Spiked Amount 50.000	Range 58 - 135		Recovery =	96.52%		
Target Compounds						Qvalue
16) Acetone	3.767	43	1324416	635.35	ug/l	99
18) Methyl Acetate	4.266	43	12647	2.04	ug/l #	85
25) 2-Butanone	6.763	43	30838	9.49	ug/l	93
37) Ethyl Acetate	6.849	43	323838	46.19	ug/l	98
51) 4-Methyl-2-Pentanone	9.936	43	10894	6.92	ug/l	91
52) Toluene	10.099	92	134495	8.63	ug/l	99

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

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