

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN061521\
 Data File : VN067359.D
 Acq On : 15 Jun 2021 20:27
 Operator : JC/MD
 Sample : M2709-17
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 SPLP-AMND-COMP-O(CL-23)

Manual Integrations
 APPROVED

MMDadoda
 6/16/2021 5:03:28 PM

Quant Time: Jun 16 02:41:27 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N061421W.M
 Quant Title : SW846 8260
 QLast Update : Tue Jun 15 05:19:42 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.088	168	329205	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.971	114	628512	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.749	117	566585	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.680	152	220508	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.440	65	315523	48.584	ug/l	0.00
Spiked Amount 50.000	Range 61 - 141		Recovery =	97.160%		
35) Dibromofluoromethane	8.024	113	163545	39.125	ug/l	0.00
Spiked Amount 50.000	Range 69 - 133		Recovery =	78.240%		
50) Toluene-d8	10.443	98	813724	50.789	ug/l	0.00
Spiked Amount 50.000	Range 65 - 126		Recovery =	101.580%		
62) 4-Bromofluorobenzene	12.736	95	311635	47.132	ug/l	0.00
Spiked Amount 50.000	Range 58 - 135		Recovery =	94.260%		
Target Compounds					Qvalue	
16) Acetone	4.299	43	24661	9.895	ug/l	96
20) Methylene Chloride	5.095	84	13229	2.551	ug/l	90
25) 2-Butanone	7.361	43	8049m	2.260	ug/l	

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

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