

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN090120\
 Data File : VN063169.D
 Acq On : 1 Sep 2020 20:53
 Operator : JC/MD
 Sample : L3851-09
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 20058

Quant Time: Sep 02 05:06:55 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N083120W.M
 Quant Title : SW846 8260
 QLast Update : Tue Sep 01 04:21:28 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.63	168	367471	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	714666	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	719346	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	291445	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	287036	47.78	ug/l	0.00
Spiked Amount 50.000			Recovery	=	95.56%	
35) Dibromofluoromethane	7.55	113	222036	48.43	ug/l	0.00
Spiked Amount 50.000			Recovery	=	96.86%	
50) Toluene-d8	10.06	98	915144	51.86	ug/l	0.00
Spiked Amount 50.000			Recovery	=	103.72%	
62) 4-Bromofluorobenzene	12.38	95	328891	53.28	ug/l	0.00
Spiked Amount 50.000			Recovery	=	106.56%	
Target Compounds						
16) Acetone	3.78	43	47536	27.837	ug/l	94
18) Methyl Acetate	4.29	43	7865m	1.796	ug/l	
20) Methylene Chloride	4.51	84	17595	3.204	ug/l	98
43) Isopropyl Acetate	8.13	43	124504	11.112	ug/l #	89

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

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