

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN090120\
 Data File : VN063172.D
 Acq On : 1 Sep 2020 22:04
 Operator : JC/MD
 Sample : L3851-18
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 RBR-78

Quant Time: Sep 02 05:15:02 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	371948	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	721683	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	723447	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	291024	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	290255	47.74	ug/l	0.00
Spiked Amount 50.000			Recovery	=	95.48%	
35) Dibromofluoromethane	7.55	113	221151	47.77	ug/l	0.00
Spiked Amount 50.000			Recovery	=	95.54%	
50) Toluene-d8	10.06	98	912872	51.23	ug/l	0.00
Spiked Amount 50.000			Recovery	=	102.46%	
62) 4-Bromofluorobenzene	12.38	95	330038	52.94	ug/l	0.00
Spiked Amount 50.000			Recovery	=	105.88%	
Target Compounds						
16) Acetone	3.78	43	33154	19.181	ug/l	92
18) Methyl Acetate	4.29	43	7462m	1.684	ug/l	
20) Methylene Chloride	4.51	84	17469	3.133	ug/l #	87
43) Isopropyl Acetate	8.13	43	113799	10.058	ug/l #	90

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

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