

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN091820\
 Data File : VN063468.D
 Acq On : 18 Sep 2020 19:25
 Operator : JC/MD
 Sample : L3871-22
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 TR-08-091720

Quant Time: Sep 21 02:31:20 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N091620W.M
 Quant Title : SW846 8260
 QLast Update : Wed Sep 16 13:05:20 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.63	168	226510	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	427961	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	434458	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	187849	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	189386	49.12	ug/l	0.00
Spiked Amount 50.000			Recovery	=	98.24%	
35) Dibromofluoromethane	7.55	113	135480	45.73	ug/l	0.00
Spiked Amount 50.000			Recovery	=	91.46%	
50) Toluene-d8	10.06	98	555565	49.87	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.74%	
62) 4-Bromofluorobenzene	12.38	95	218659	53.54	ug/l	0.00
Spiked Amount 50.000			Recovery	=	107.08%	
Target Compounds						
16) Acetone	3.78	43	35725	28.829	ug/l	94
18) Methyl Acetate	4.28	43	3338	1.061	ug/l	91
20) Methylene Chloride	4.51	84	14787	4.710	ug/l	88
43) Isopropyl Acetate	8.13	43	84649	10.579	ug/l	98

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

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