

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN092220\
 Data File : VN063571.D
 Acq On : 23 Sep 2020 7:40
 Operator : JC/MD
 Sample : L3866-10
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 50 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 OK-01-092120

Manual Integrations
 APPROVED

MMDadoda
 9/23/2020 1:06:54 PM

Quant Time: Sep 23 08:42:05 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	200919	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	389962	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	391506	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	159224	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	175626	51.35	ug/l	0.00
Spiked Amount 50.000			Recovery	=	102.70%	
35) Dibromofluoromethane	7.55	113	124792	46.23	ug/l	0.00
Spiked Amount 50.000			Recovery	=	92.46%	
50) Toluene-d8	10.06	98	498038	49.06	ug/l	0.00
Spiked Amount 50.000			Recovery	=	98.12%	
62) 4-Bromofluorobenzene	12.38	95	191793	51.54	ug/l	0.00
Spiked Amount 50.000			Recovery	=	103.08%	
Target Compounds						
16) Acetone	3.78	43	28193	25.649	ug/l	99
18) Methyl Acetate	4.30	43	4561m	1.635	ug/l	
20) Methylene Chloride	4.50	84	12451	4.471	ug/l	88
43) Isopropyl Acetate	8.13	43	58854	8.072	ug/l	100

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

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