

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN092420\
 Data File : VN063661.D
 Acq On : 25 Sep 2020 5:25
 Operator : JC/MD
 Sample : L4106-02
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 47 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 GLOUCESTER-SP

Manual Integrations
 APPROVED

MMDadoda
 9/28/2020 5:15:27 PM

Quant Time: Sep 25 06:22:20 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	198591	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	392337	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	396687	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	161514	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	175557	51.93	ug/l	0.00
Spiked Amount 50.000			Recovery	=	103.86%	
35) Dibromofluoromethane	7.55	113	126316	46.51	ug/l	0.00
Spiked Amount 50.000			Recovery	=	93.02%	
50) Toluene-d8	10.06	98	501724	49.13	ug/l	0.00
Spiked Amount 50.000			Recovery	=	98.26%	
62) 4-Bromofluorobenzene	12.38	95	192946	51.54	ug/l	0.00
Spiked Amount 50.000			Recovery	=	103.08%	
Target Compounds						
16) Acetone	3.78	43	15231	14.019	ug/l	98
18) Methyl Acetate	4.28	43	4841m	1.755	ug/l	
20) Methylene Chloride	4.50	84	16067	5.837	ug/l	93
43) Isopropyl Acetate	8.13	43	71498	9.747	ug/l	99

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

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