

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN090420\
 Data File : VN063248.D
 Acq On : 4 Sep 2020 15:25
 Operator : JC/MD
 Sample : L3902-18
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 SA-PZ172I-20200902

Manual Integrations
 APPROVED

MMDadoda
 9/8/2020 1:33:43 PM

Quant Time: Sep 07 04:05:22 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	273728	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	529610	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	523441	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	202516	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	218901	48.92	ug/l	0.00
Spiked Amount 50.000			Recovery	=	97.84%	
35) Dibromofluoromethane	7.55	113	168315	49.54	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.08%	
50) Toluene-d8	10.06	98	672942	51.46	ug/l	0.00
Spiked Amount 50.000			Recovery	=	102.92%	
62) 4-Bromofluorobenzene	12.38	95	228452	49.94	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.88%	
Target Compounds						
3) Chloromethane	2.03	50	5739	1.462	ug/l #	83
12) 1,1-Dichloroethene	3.69	96	2070m	0.689	ug/l	
19) Methyl tert-butyl Ether	5.00	73	10546m	0.947	ug/l	
20) Methylene Chloride	4.50	84	5301m	1.000	ug/l	
24) 1,1-Dichloroethane	5.80	63	44672	6.780	ug/l	98

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

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