

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN092320\
 Data File : VN063600.D
 Acq On : 23 Sep 2020 19:42
 Operator : JC/MD
 Sample : L4083-06
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 MSVOA_N
ClientSampleId :
 RE135D2-20200917

Manual Integrations
APPROVED
 MMDadoda
 9/24/2020 1:38:08 PM

Quant Time: Sep 24 04:09:22 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	194937	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	375946	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	374069	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	154464	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	174770	52.67	ug/l	0.00
Spiked Amount	50.000		Recovery	=	105.34%	
35) Dibromofluoromethane	7.55	113	125352	48.16	ug/l	0.00
Spiked Amount	50.000		Recovery	=	96.32%	
50) Toluene-d8	10.06	98	480626	49.11	ug/l	0.00
Spiked Amount	50.000		Recovery	=	98.22%	
62) 4-Bromofluorobenzene	12.38	95	185369	51.67	ug/l	0.00
Spiked Amount	50.000		Recovery	=	103.34%	
Target Compounds						
3) Chloromethane	2.04	50	2735	0.936	ug/l	89
16) Acetone	3.79	43	5569m	5.222	ug/l	
44) Trichloroethene	8.80	130	2443	0.881	ug/l	99
45) 1,2-Dichloropropane	9.09	63	6432	2.066	ug/l	99
52) Toluene	10.13	92	3478	0.506	ug/l	83

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

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