

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN100120\
 Data File : VN063799.D
 Acq On : 2 Oct 2020 3:20
 Operator : JC/MD
 Sample : L4190-02
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 42 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PAV-B14-092920-0-1

Manual Integrations
 APPROVED

MMDadoda
 10/2/2020 3:12:10 PM

Quant Time: Oct 02 04:46:23 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N093020W.M
 Quant Title : SW846 8260
 QLast Update : Thu Oct 01 01:46:38 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.63	168	214584	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	416765	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	419399	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	171560	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	193881	54.69	ug/l	0.00
Spiked Amount 50.000			Recovery	=	109.38%	
35) Dibromofluoromethane	7.55	113	137131	47.42	ug/l	0.00
Spiked Amount 50.000			Recovery	=	94.84%	
50) Toluene-d8	10.06	98	525458	47.93	ug/l	0.00
Spiked Amount 50.000			Recovery	=	95.86%	
62) 4-Bromofluorobenzene	12.38	95	202561	49.15	ug/l	0.00
Spiked Amount 50.000			Recovery	=	98.30%	
Target Compounds						
16) Acetone	3.78	43	32911	30.158	ug/l	99
18) Methyl Acetate	4.29	43	4937m	1.742	ug/l	
20) Methylene Chloride	4.51	84	8866	3.638	ug/l	96
43) Isopropyl Acetate	8.13	43	75939	10.975	ug/l	100

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

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