

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN100120\
 Data File : VN063800.D
 Acq On : 2 Oct 2020 3:44
 Operator : JC/MD
 Sample : L4190-08
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 43 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PAV-B14-092920-15-16

Manual Integrations
 APPROVED

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

Quant Time: Oct 02 04:49:01 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	7.62	168	214829	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	403608	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	419385	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	170896	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	189997	53.53	ug/l	0.00
Spiked Amount 50.000			Recovery	=	107.06%	
35) Dibromofluoromethane	7.55	113	136563	48.76	ug/l	0.00
Spiked Amount 50.000			Recovery	=	97.52%	
50) Toluene-d8	10.06	98	522509	49.22	ug/l	0.00
Spiked Amount 50.000			Recovery	=	98.44%	
62) 4-Bromofluorobenzene	12.38	95	205555	51.51	ug/l	0.00
Spiked Amount 50.000			Recovery	=	103.02%	
Target Compounds						
16) Acetone	3.78	43	35765	32.736	ug/l	95
18) Methyl Acetate	4.29	43	4386	1.546	ug/l #	77
20) Methylene Chloride	4.51	84	7733m	3.182	ug/l	
43) Isopropyl Acetate	8.13	43	73697	10.998	ug/l	98

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

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