

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN100120\
 Data File : VN063801.D
 Acq On : 2 Oct 2020 4:08
 Operator : JC/MD
 Sample : L4190-10
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 44 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PAV-B18-092920-0-1

Manual Integrations
 APPROVED

MMDadoda
 10/2/2020 3:12:14 PM

Quant Time: Oct 02 05:48:55 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	209547	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	395998	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	409497	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	170129	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	189430	54.72	ug/l	0.00
Spiked Amount 50.000			Recovery	=	109.44%	
35) Dibromofluoromethane	7.55	113	135096	49.17	ug/l	0.00
Spiked Amount 50.000			Recovery	=	98.34%	
50) Toluene-d8	10.06	98	512583	49.21	ug/l	0.00
Spiked Amount 50.000			Recovery	=	98.42%	
62) 4-Bromofluorobenzene	12.38	95	203173	51.89	ug/l	0.00
Spiked Amount 50.000			Recovery	=	103.78%	
Target Compounds						
16) Acetone	3.78	43	22935	21.522	ug/l	96
18) Methyl Acetate	4.30	43	4357m	1.575	ug/l	
20) Methylene Chloride	4.51	84	9047	3.797	ug/l	97
43) Isopropyl Acetate	8.13	43	74854	11.386	ug/l	98

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

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