

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN100620\
 Data File : VN063923.D
 Acq On : 7 Oct 2020 4:17
 Operator : JC/MD
 Sample : L4242-18
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 42 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PAV-B24-100120-DP

Manual Integrations
 APPROVED

MMDadoda
 10/7/2020 12:12:18 PM

Quant Time: Oct 07 06:03:28 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	229139	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	412844	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	437560	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	173658	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	196262	51.85	ug/l	0.00
Spiked Amount 50.000			Recovery	=	103.70%	
35) Dibromofluoromethane	7.55	113	143777	50.19	ug/l	0.00
Spiked Amount 50.000			Recovery	=	100.38%	
50) Toluene-d8	10.06	98	543766	50.07	ug/l	0.00
Spiked Amount 50.000			Recovery	=	100.14%	
62) 4-Bromofluorobenzene	12.38	95	210704	51.62	ug/l	0.00
Spiked Amount 50.000			Recovery	=	103.24%	
Target Compounds						
16) Acetone	3.78	43	15857	13.608	ug/l	98
18) Methyl Acetate	4.28	43	4058m	1.341	ug/l	
20) Methylene Chloride	4.50	84	16023m	6.089	ug/l	
40) Benzene	8.01	78	18024	1.509	ug/l	96
43) Isopropyl Acetate	8.13	43	81908	11.950	ug/l	99

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

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