

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN100620\
 Data File : VN063925.D
 Acq On : 7 Oct 2020 5:06
 Operator : JC/MD
 Sample : L4247-13
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 44 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 B-19(4-5)

Manual Integrations
 APPROVED

MMDadoda
 10/7/2020 12:12:20 PM

Quant Time: Oct 07 06:09:36 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	227651	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	422684	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	442315	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	173145	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	196676	52.29	ug/l	0.00
Spiked Amount 50.000			Recovery	=	104.58%	
35) Dibromofluoromethane	7.55	113	145275	49.53	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.06%	
50) Toluene-d8	10.06	98	554136	49.84	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.68%	
62) 4-Bromofluorobenzene	12.38	95	209533	50.13	ug/l	0.00
Spiked Amount 50.000			Recovery	=	100.26%	
Target Compounds						
16) Acetone	3.78	43	26656	23.024	ug/l	99
17) Carbon Disulfide	4.02	76	2748	1.138	ug/l #	85
18) Methyl Acetate	4.29	43	4521m	1.504	ug/l	
20) Methylene Chloride	4.50	84	9117	3.529	ug/l	96
25) 2-Butanone	6.79	43	8730m	5.319	ug/l	
43) Isopropyl Acetate	8.13	43	87457	12.463	ug/l	98

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

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