

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN101220\
 Data File : VN064088.D
 Acq On : 12 Oct 2020 19:26
 Operator : JC/MD
 Sample : L4333-04
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 OH-WC-J6-3-6

Manual Integrations
 APPROVED

MMDadoda
 10/13/2020 4:09:52 PM

Quant Time: Oct 13 02:34:28 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	223073	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	410038	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	434112	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	171070	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	188223	51.07	ug/l	0.00
Spiked Amount 50.000			Recovery	=	102.14%	
35) Dibromofluoromethane	7.55	113	144449	50.77	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.54%	
50) Toluene-d8	10.06	98	545287	50.56	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.12%	
62) 4-Bromofluorobenzene	12.38	95	213044	52.55	ug/l	0.00
Spiked Amount 50.000			Recovery	=	105.10%	
Target Compounds						
16) Acetone	3.78	43	31160	27.467	ug/l	96
17) Carbon Disulfide	4.02	76	3073m	1.200	ug/l	
18) Methyl Acetate	4.29	43	9833	3.338	ug/l #	70
20) Methylene Chloride	4.51	84	10621	4.177	ug/l	95
43) Isopropyl Acetate	8.13	43	57352	8.425	ug/l	99

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

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