

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN111020\
 Data File : VN064526.D
 Acq On : 10 Nov 2020 15:24
 Operator : JC/MD
 Sample : L4684-03
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 BP-VPB178-GW-398-400

Manual Integrations
 APPROVED

MMDadoda
 11/11/2020 1:01:46 PM

Quant Time: Nov 11 07:25:50 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N110420W.M
 Quant Title : SW846 8260
 QLast Update : Wed Nov 04 12:09:03 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.63	168	242411	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	490699	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	515517	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	183520	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	237313	56.20	ug/l	0.00
Spiked Amount 50.000			Recovery	=	112.40%	
35) Dibromofluoromethane	7.55	113	174386	52.20	ug/l	0.00
Spiked Amount 50.000			Recovery	=	104.40%	
50) Toluene-d8	10.06	98	645795	50.01	ug/l	0.00
Spiked Amount 50.000			Recovery	=	100.02%	
62) 4-Bromofluorobenzene	12.38	95	239845	49.11	ug/l	0.00
Spiked Amount 50.000			Recovery	=	98.22%	
Target Compounds						
12) 1,1-Dichloroethene	3.69	96	1799m	0.694	ug/l	Qvalue
16) Acetone	3.79	43	8183	5.088	ug/l	94
19) Methyl tert-butyl Ether	5.01	73	4187	0.449	ug/l	97
27) cis-1,2-Dichloroethene	6.79	96	21483	6.298	ug/l	88
30) Chloroform	7.33	83	1902	0.306	ug/l	88
44) Trichloroethene	8.80	130	221112	57.532	ug/l	98
45) 1,2-Dichloropropane	9.09	63	22027	5.419	ug/l	98
64) Tetrachloroethene	10.60	164	14932	3.760	ug/l	94

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

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