

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN040720\
 Data File : VN060884.D
 Acq On : 7 Apr 2020 17:57
 Operator : JC/MD
 Sample : L2178-05
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 MW-1A

Manual Integrations
 APPROVED

MMDadoda
 4/8/2020 9:39:03 AM

Quant Time: Apr 08 08:33:08 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N031820W.M
 Quant Title : SW846 8260
 QLast Update : Wed Mar 18 08:49:09 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.64	168	196480	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.57	114	325476	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.40	117	300579	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.34	152	137643	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.00	65	124937	46.32	ug/l	0.00
Spiked Amount 50.000			Recovery	=	92.64%	
35) Dibromofluoromethane	7.57	113	96643	49.11	ug/l	0.00
Spiked Amount 50.000			Recovery	=	98.22%	
50) Toluene-d8	10.08	98	400432	49.81	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.62%	
62) 4-Bromofluorobenzene	12.40	95	148187	50.30	ug/l	0.00
Spiked Amount 50.000			Recovery	=	100.60%	
Target Compounds						
					Qvalue	
3) Chloromethane	2.04	50	1124	0.317	ug/l	96
27) cis-1,2-Dichloroethene	6.81	96	6057	2.351	ug/l	86
30) Chloroform	7.34	83	3556	0.861	ug/l	93
40) Benzene	8.02	78	7884	0.829	ug/l	96
67) Ethyl Benzene	11.50	91	14566	1.314	ug/l	96
73) Isopropylbenzene	12.25	105	18991	1.839	ug/l	99
78) n-propylbenzene	12.59	91	18580	1.522	ug/l	100
84) 1,2,4-Trimethylbenzene	13.04	105	4321	0.498	ug/l	97
85) sec-Butylbenzene	13.17	105	8666	0.887	ug/l	85
89) n-Butylbenzene	13.61	91	5430m	0.684	ug/l	
95) Naphthalene	15.13	128	13239	1.708	ug/l #	96

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

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