

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN111120\
 Data File : VN064551.D
 Acq On : 11 Nov 2020 13:36
 Operator : JC/MD
 Sample : L4684-02DL 4X
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 BP-VPB178-DUP-20201106DL

Manual Integrations
 APPROVED

MMDadoda
 11/12/2020 9:12:27 AM

Quant Time: Nov 12 02:12:34 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	239247	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	491040	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	498258	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	178022	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	238677	57.27	ug/l	0.00
Spiked Amount 50.000			Recovery	=	114.54%	
35) Dibromofluoromethane	7.55	113	174573	52.22	ug/l	0.00
Spiked Amount 50.000			Recovery	=	104.44%	
50) Toluene-d8	10.06	98	628531	48.64	ug/l	0.00
Spiked Amount 50.000			Recovery	=	97.28%	
62) 4-Bromofluorobenzene	12.38	95	230953	47.26	ug/l	0.00
Spiked Amount 50.000			Recovery	=	94.52%	
Target Compounds						
					Qvalue	
11) Tert butyl alcohol	4.75	59	7498	16.315	ug/l #	71
12) 1,1-Dichloroethene	3.69	96	4880	1.907	ug/l #	70
22) Diisopropyl ether	5.92	45	2506m	0.236	ug/l	
24) 1,1-Dichloroethane	5.80	63	35954	6.158	ug/l	97
40) Benzene	8.01	78	836934	54.904	ug/l	99
42) 1,2-Dichloroethane	8.09	62	30344	4.933	ug/l	99
44) Trichloroethene	8.80	130	2567	0.667	ug/l	76
69) o-Xylene	11.92	106	6869	1.009	ug/l	94
73) Isopropylbenzene	12.23	105	18366	1.298	ug/l	97

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

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