

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN111020\
 Data File : VN064533.D
 Acq On : 10 Nov 2020 18:13
 Operator : JC/MD
 Sample : L4684-08
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 BP-VPB178-GW-498-500

Manual Integrations
 APPROVED

MMDadoda
 11/11/2020 1:01:57 PM

Quant Time: Nov 11 07:50:08 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	241706	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	501035	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	511109	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	190217	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	234495	55.70	ug/l	0.00
Spiked Amount 50.000			Recovery	=	111.40%	
35) Dibromofluoromethane	7.55	113	172494	50.56	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.12%	
50) Toluene-d8	10.06	98	633038	48.01	ug/l	0.00
Spiked Amount 50.000			Recovery	=	96.02%	
62) 4-Bromofluorobenzene	12.38	95	251954	50.53	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.06%	
Target Compounds						
					Qvalue	
11) Tert butyl alcohol	4.76	59	13548	29.180	ug/l #	71
12) 1,1-Dichloroethene	3.69	96	82425	31.886	ug/l	98
20) Methylene Chloride	4.50	84	12876	4.047	ug/l	86
22) Diisopropyl ether	5.92	45	14395m	1.341	ug/l	
24) 1,1-Dichloroethane	5.80	63	251315	42.607	ug/l	100
32) 1,1,1-Trichloroethane	7.54	97	30702	5.583	ug/l	99
40) Benzene	8.00	78	2539604	163.279	ug/l	99
42) 1,2-Dichloroethane	8.09	62	203933	32.490	ug/l	98
44) Trichloroethene	8.80	130	4436	1.130	ug/l	94
49) 1,4-Dioxane	9.17	88	4551	79.826	ug/l #	93
64) Tetrachloroethene	10.60	164	3331	0.846	ug/l #	92
69) o-Xylene	11.92	106	62245	8.912	ug/l	93
73) Isopropylbenzene	12.23	105	55969	3.702	ug/l	99

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

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