

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN111020\
 Data File : VN064530.D
 Acq On : 10 Nov 2020 17:00
 Operator : JC/MD
 Sample : L4684-07
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 BP-VPB178-GW-478-480

Manual Integrations
 APPROVED

MMDadoda
 11/11/2020 1:01:53 PM

Quant Time: Nov 11 07:44:00 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	237725	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	486628	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	500267	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	190539	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	229424	55.41	ug/l	0.00
Spiked Amount 50.000			Recovery	=	110.82%	
35) Dibromofluoromethane	7.55	113	170220	51.38	ug/l	0.00
Spiked Amount 50.000			Recovery	=	102.76%	
50) Toluene-d8	10.06	98	619860	48.40	ug/l	0.00
Spiked Amount 50.000			Recovery	=	96.80%	
62) 4-Bromofluorobenzene	12.38	95	244009	50.38	ug/l	0.00
Spiked Amount 50.000			Recovery	=	100.76%	
Target Compounds						
					Qvalue	
11) Tert butyl alcohol	4.75	59	12670	27.746	ug/l #	71
12) 1,1-Dichloroethene	3.70	96	21435	8.431	ug/l	93
22) Diisopropyl ether	5.90	45	9746m	0.923	ug/l	
24) 1,1-Dichloroethane	5.80	63	160629	27.688	ug/l	99
32) 1,1,1-Trichloroethane	7.53	97	8548	1.580	ug/l	97
40) Benzene	8.01	78	1788693	118.406	ug/l	100
42) 1,2-Dichloroethane	8.09	62	133890	21.962	ug/l	98
44) Trichloroethene	8.80	130	12701	3.332	ug/l	99
49) 1,4-Dioxane	9.17	88	1195	21.581	ug/l #	74
52) Toluene	10.13	92	8305	0.901	ug/l	89
64) Tetrachloroethene	10.60	164	3540	0.919	ug/l	88
68) m/p-Xylenes	11.59	106	2621	0.370	ug/l	92
69) o-Xylene	11.92	106	49859	7.293	ug/l	99
73) Isopropylbenzene	12.23	105	39462	2.606	ug/l	95

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

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