

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN111020\
 Data File : VN064531.D
 Acq On : 10 Nov 2020 17:25
 Operator : JC/MD
 Sample : L4684-11
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 BP-VPB178-GW-518-520

Manual Integrations
 APPROVED

MMDadoda
 11/11/2020 1:01:55 PM

Quant Time: Nov 11 04:07:16 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	237170	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	491632	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	504464	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	189285	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	231991	56.16	ug/l	0.00
Spiked Amount 50.000			Recovery	=	112.32%	
35) Dibromofluoromethane	7.55	113	169602	50.67	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.34%	
50) Toluene-d8	10.06	98	624910	48.30	ug/l	0.00
Spiked Amount 50.000			Recovery	=	96.60%	
62) 4-Bromofluorobenzene	12.38	95	248162	50.72	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.44%	
Target Compounds						
					Qvalue	
12) 1,1-Dichloroethene	3.69	96	68444	26.984	ug/l	97
16) Acetone	3.77	43	15639	9.938	ug/l #	80
20) Methylene Chloride	4.51	84	9452	3.027	ug/l	94
22) Diisopropyl ether	5.92	45	10570m	1.004	ug/l	
24) 1,1-Dichloroethane	5.80	63	195683	33.810	ug/l	100
32) 1,1,1-Trichloroethane	7.53	97	30431	5.640	ug/l	97
40) Benzene	8.01	78	1458446	95.562	ug/l	100
42) 1,2-Dichloroethane	8.09	62	161827	26.275	ug/l	98
44) Trichloroethene	8.80	130	4586	1.191	ug/l	82
49) 1,4-Dioxane	9.17	88	3814	68.178	ug/l #	77
64) Tetrachloroethene	10.60	164	2998	0.772	ug/l	93
69) o-Xylene	11.92	106	102155	14.819	ug/l	100
73) Isopropylbenzene	12.23	105	28684	1.907	ug/l	99

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

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