

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN111620\
 Data File : VN064661.D
 Acq On : 16 Nov 2020 15:00
 Operator : JC/MD
 Sample : L4774-01
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 MW-1

Manual Integrations
 APPROVED

MMDadoda
 11/17/2020 7:02:13 PM

Quant Time: Nov 17 04:39:54 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	7.62	168	226308	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	453188	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	489310	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	217911	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	220131	55.84	ug/l	0.00
Spiked Amount 50.000			Recovery	=	111.68%	
35) Dibromofluoromethane	7.55	113	163817	53.09	ug/l	0.00
Spiked Amount 50.000			Recovery	=	106.18%	
50) Toluene-d8	10.06	98	600450	50.35	ug/l	0.00
Spiked Amount 50.000			Recovery	=	100.70%	
62) 4-Bromofluorobenzene	12.38	95	247662	54.91	ug/l	0.00
Spiked Amount 50.000			Recovery	=	109.82%	
Target Compounds						
						Qvalue
16) Acetone	3.80	43	7832m	5.216	ug/l	
39) Methylcyclohexane	9.05	83	6225m	1.450	ug/l	
52) Toluene	10.12	92	35356	4.120	ug/l	97
67) Ethyl Benzene	11.48	91	327414	17.160	ug/l	100
68) m/p-Xylenes	11.60	106	159203	23.008	ug/l	99
69) o-Xylene	11.92	106	35675	5.335	ug/l	92
73) Isopropylbenzene	12.22	105	11937	0.689	ug/l	97
78) n-propylbenzene	12.57	91	9675	0.490	ug/l #	54
80) 1,3,5-Trimethylbenzene	12.71	105	539691	37.770	ug/l	98
84) 1,2,4-Trimethylbenzene	13.02	105	283876	19.763	ug/l	99
85) sec-Butylbenzene	13.15	105	3959m	0.280	ug/l	
86) p-Isopropyltoluene	13.26	119	20387	1.604	ug/l	97
89) n-Butylbenzene	13.56	91	75400	6.898	ug/l #	1

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

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