

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN111920\
 Data File : VN064768.D
 Acq On : 20 Nov 2020 8:19
 Operator : JC/MD
 Sample : L4779-02 50X
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 50 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 1641

Manual Integrations
 APPROVED

MMDadoda
 11/24/2020 3:11:41 PM

Quant Time: Nov 20 14:52:00 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N111820W.M
 Quant Title : SW846 8260
 QLast Update : Wed Nov 18 15:29:54 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.63	168	217072	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	523885	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	502369	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	288322	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	254967	66.77	ug/l	0.00
Spiked Amount 50.000			Recovery	=	133.54%	
35) Dibromofluoromethane	7.55	113	152430	51.12	ug/l	0.00
Spiked Amount 50.000			Recovery	=	102.24%	
50) Toluene-d8	10.06	98	569194	42.00	ug/l	0.00
Spiked Amount 50.000			Recovery	=	84.00%	
62) 4-Bromofluorobenzene	12.37	95	315058	62.20	ug/l	0.00
Spiked Amount 50.000			Recovery	=	124.40%	
Target Compounds						
16) Acetone	3.78	43	8412	13.985	ug/l	98
30) Chloroform	7.33	83	8696	1.614	ug/l	91
44) Trichloroethene	8.81	130	2717m	0.630	ug/l	

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

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