

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN100820\
 Data File : VN063993.D
 Acq On : 8 Oct 2020 16:06
 Operator : JC/MD
 Sample : L4255-04
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 SA-PZ124-20200930

Manual Integrations
 APPROVED

MMDadoda
 10/9/2020 4:32:17 PM

Quant Time: Oct 09 13:31:40 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N093020W.M
 Quant Title : SW846 8260
 QLast Update : Thu Oct 01 01:46:38 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.63	168	245433	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	431581	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	447088	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	172048	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	203097	50.09	ug/l	0.00
Spiked Amount 50.000			Recovery	=	100.18%	
35) Dibromofluoromethane	7.55	113	153783	51.35	ug/l	0.00
Spiked Amount 50.000			Recovery	=	102.70%	
50) Toluene-d8	10.06	98	557269	49.09	ug/l	0.00
Spiked Amount 50.000			Recovery	=	98.18%	
62) 4-Bromofluorobenzene	12.38	95	208787	48.93	ug/l	0.00
Spiked Amount 50.000			Recovery	=	97.86%	
Target Compounds						
6) Chloroethane	2.68	64	5178	2.841	ug/l	97
12) 1,1-Dichloroethene	3.70	96	3975	1.608	ug/l #	72
16) Acetone	3.78	43	5912m	4.737	ug/l	
24) 1,1-Dichloroethane	5.79	63	56623	10.483	ug/l	99

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

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