

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN081920\
 Data File : VN063011.D
 Acq On : 19 Aug 2020 14:15
 Operator : JC/MD
 Sample : L3718-04
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 BP-VPB181-DUP-20200812

Manual Integrations
 APPROVED

MMDadoda
 8/20/2020 10:09:36 AM

Quant Time: Aug 19 15:37:17 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N081820W.M
 Quant Title : SW846 8260
 QLast Update : Wed Aug 19 06:09:27 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.63	168	147702	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	244651	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	253171	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	92762	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	92461	52.82	ug/l	0.00
Spiked Amount 50.000			Recovery	=	105.64%	
35) Dibromofluoromethane	7.55	113	82108	51.51	ug/l	0.00
Spiked Amount 50.000			Recovery	=	103.02%	
50) Toluene-d8	10.06	98	314599	51.38	ug/l	0.00
Spiked Amount 50.000			Recovery	=	102.76%	
62) 4-Bromofluorobenzene	12.38	95	105356	48.76	ug/l	0.00
Spiked Amount 50.000			Recovery	=	97.52%	
Target Compounds						
12) 1,1-Dichloroethene	3.70	96	2366	1.887	ug/l	91
30) Chloroform	7.33	83	2462m	0.872	ug/l	
44) Trichloroethene	8.80	130	6417	3.773	ug/l	95
64) Tetrachloroethene	10.61	164	951	0.571	ug/l	90

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

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