

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN090920\
 Data File : VN063283.D
 Acq On : 9 Sep 2020 15:28
 Operator : JC/MD
 Sample : L3931-04
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 SP-1-2

Quant Time: Sep 10 04:27:00 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N083120W.M
 Quant Title : SW846 8260
 QLast Update : Tue Sep 01 04:21:28 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.63	168	417615	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	799949	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	780054	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	318634	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	330197	48.37	ug/l	0.00
Spiked Amount 50.000			Recovery	=	96.74%	
35) Dibromofluoromethane	7.55	113	251556	49.02	ug/l	0.00
Spiked Amount 50.000			Recovery	=	98.04%	
50) Toluene-d8	10.06	98	1016466	51.46	ug/l	0.00
Spiked Amount 50.000			Recovery	=	102.92%	
62) 4-Bromofluorobenzene	12.38	95	391885	56.71	ug/l	0.00
Spiked Amount 50.000			Recovery	=	113.42%	
Target Compounds						
16) Acetone	3.78	43	36589	18.854	ug/l	100
18) Methyl Acetate	4.29	43	6831	1.373	ug/l #	88
20) Methylene Chloride	4.51	84	20682	3.331	ug/l	96
43) Isopropyl Acetate	8.13	43	148586	11.847	ug/l #	90

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

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