

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN090920\
 Data File : VN063282.D
 Acq On : 9 Sep 2020 15:04
 Operator : JC/MD
 Sample : L3931-02
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 SP-1-1

Quant Time: Sep 10 04:24:55 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	425149	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	791188	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	773141	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	329938	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	325723	46.87	ug/l	0.00
Spiked Amount 50.000			Recovery	=	93.74%	
35) Dibromofluoromethane	7.55	113	247344	48.73	ug/l	0.00
Spiked Amount 50.000			Recovery	=	97.46%	
50) Toluene-d8	10.06	98	1013707	51.89	ug/l	0.00
Spiked Amount 50.000			Recovery	=	103.78%	
62) 4-Bromofluorobenzene	12.38	95	393783	57.62	ug/l	0.00
Spiked Amount 50.000			Recovery	=	115.24%	
Target Compounds						
16) Acetone	3.78	43	38410	19.442	ug/l	91
18) Methyl Acetate	4.28	43	5240	1.034	ug/l #	56
20) Methylene Chloride	4.51	84	20244	3.184	ug/l	93
43) Isopropyl Acetate	8.13	43	139899	11.278	ug/l #	90

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

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