

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN031620\
 Data File : VN060523.D
 Acq On : 16 Mar 2020 15:14
 Operator : JC/MD
 Sample : L1933-06
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 72-11976

Quant Time: Mar 17 09:24:13 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N022720W.M
 Quant Title : SW846 8260
 QLast Update : Thu Feb 27 13:52:50 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.64	168	213267	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.57	114	357051	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.40	117	333425	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.34	152	149133	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.00	65	138653	50.67	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.34%	
35) Dibromofluoromethane	7.57	113	104265	62.91	ug/l	0.00
Spiked Amount 50.000			Recovery	=	125.82%	
50) Toluene-d8	10.08	98	443222	52.84	ug/l	0.00
Spiked Amount 50.000			Recovery	=	105.68%	
62) 4-Bromofluorobenzene	12.40	95	163540	51.43	ug/l	0.00
Spiked Amount 50.000			Recovery	=	102.86%	
Target Compounds						
16) Acetone	3.79	43	9894	11.197	ug/l	100
20) Methylene Chloride	4.53	84	5906	2.434	ug/l	93
43) Isopropyl Acetate	8.14	43	62959	11.161	ug/l #	90
51) 4-Methyl-2-Pentanone	9.97	43	5188	1.622	ug/l	95

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

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