

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN031320\
 Data File : VN060510.D
 Acq On : 13 Mar 2020 18:05
 Operator : JC/MD
 Sample : L1907-04
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 TP-6-SA

Quant Time: Mar 16 09:18:59 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	212677	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.57	114	356004	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.40	117	340687	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.34	152	158938	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.00	65	139094	50.97	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.94%	
35) Dibromofluoromethane	7.56	113	104665	63.34	ug/l	0.00
Spiked Amount 50.000			Recovery	=	126.68%	
50) Toluene-d8	10.08	98	440575	52.68	ug/l	0.00
Spiked Amount 50.000			Recovery	=	105.36%	
62) 4-Bromofluorobenzene	12.40	95	166622	52.56	ug/l	0.00
Spiked Amount 50.000			Recovery	=	105.12%	
Target Compounds						
16) Acetone	3.79	43	7729	8.771	ug/l	92
20) Methylene Chloride	4.52	84	5538	2.288	ug/l	94
43) Isopropyl Acetate	8.14	43	63457	11.283	ug/l #	90
51) 4-Methyl-2-Pentanone	9.97	43	5113	1.603	ug/l	97

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

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