

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN031620\
 Data File : VN060524.D
 Acq On : 16 Mar 2020 15:37
 Operator : JC/MD
 Sample : L1942-04
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 TP-10-SA

Quant Time: Mar 17 09:25:49 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	215976	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.57	114	358870	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.40	117	338262	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.34	152	154866	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.00	65	140581	50.73	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.46%	
35) Dibromofluoromethane	7.56	113	105268	63.20	ug/l	0.00
Spiked Amount 50.000			Recovery	=	126.40%	
50) Toluene-d8	10.08	98	442858	52.53	ug/l	0.00
Spiked Amount 50.000			Recovery	=	105.06%	
62) 4-Bromofluorobenzene	12.40	95	166054	51.96	ug/l	0.00
Spiked Amount 50.000			Recovery	=	103.92%	
Target Compounds						
16) Acetone	3.78	43	7226	8.075	ug/l	92
20) Methylene Chloride	4.52	84	6945	2.826	ug/l	98
43) Isopropyl Acetate	8.14	43	63586	11.215	ug/l #	89
51) 4-Methyl-2-Pentanone	9.97	43	4461	1.388	ug/l	97

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

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