

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN090120\
 Data File : VN063163.D
 Acq On : 1 Sep 2020 18:29
 Operator : JC/MD
 Sample : L3849-06
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 232

Quant Time: Sep 02 04:48:51 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	379301	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	729072	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	726802	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	288768	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	291626	47.03	ug/l	0.00
Spiked Amount 50.000			Recovery	=	94.06%	
35) Dibromofluoromethane	7.55	113	231015	49.39	ug/l	0.00
Spiked Amount 50.000			Recovery	=	98.78%	
50) Toluene-d8	10.06	98	935510	51.96	ug/l	0.00
Spiked Amount 50.000			Recovery	=	103.92%	
62) 4-Bromofluorobenzene	12.38	95	333313	52.93	ug/l	0.00
Spiked Amount 50.000			Recovery	=	105.86%	
Target Compounds						
16) Acetone	3.78	43	37337	21.183	ug/l	91
18) Methyl Acetate	4.28	43	8298	1.836	ug/l #	73
20) Methylene Chloride	4.51	84	16024	2.768	ug/l	94
43) Isopropyl Acetate	8.13	43	137517	12.031	ug/l #	90

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

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