

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN063020\
 Data File : VN062279.D
 Acq On : 1 Jul 2020 8:20
 Operator : JC/MD
 Sample : L3153-38
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 54 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 SP2-J

Quant Time: Jul 01 09:56:44 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N062420W.M
 Quant Title : SW846 8260
 QLast Update : Wed Jun 24 15:33:00 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.63	168	193151	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	410330	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	398924	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	144955	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	155277	57.29	ug/l	0.00
Spiked Amount 50.000			Recovery	=	114.58%	
35) Dibromofluoromethane	7.55	113	124975	47.47	ug/l	0.00
Spiked Amount 50.000			Recovery	=	94.94%	
50) Toluene-d8	10.06	98	529154	50.29	ug/l	0.00
Spiked Amount 50.000			Recovery	=	100.58%	
62) 4-Bromofluorobenzene	12.38	95	172355	48.48	ug/l	0.00
Spiked Amount 50.000			Recovery	=	96.96%	
Target Compounds						
16) Acetone	3.78	43	49432	54.939	ug/l	99
18) Methyl Acetate	4.29	43	7271	3.374	ug/l #	70
20) Methylene Chloride	4.50	84	18338	7.184	ug/l	89
43) Isopropyl Acetate	8.13	43	43708	7.851	ug/l	99

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

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