

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN063020\
 Data File : VN062268.D
 Acq On : 1 Jul 2020 3:54
 Operator : JC/MD
 Sample : L3153-05
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 43 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 SP1-B

Manual Integrations
 APPROVED

MMDadoda
 7/1/2020 11:04:07 AM

Quant Time: Jul 01 06:07:37 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	225326	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	442998	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	433431	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	161171	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	168080	53.15	ug/l	0.00
Spiked Amount 50.000			Recovery	=	106.30%	
35) Dibromofluoromethane	7.55	113	136916	48.18	ug/l	0.00
Spiked Amount 50.000			Recovery	=	96.36%	
50) Toluene-d8	10.06	98	575364	50.65	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.30%	
62) 4-Bromofluorobenzene	12.38	95	192015	50.03	ug/l	0.00
Spiked Amount 50.000			Recovery	=	100.06%	
Target Compounds						
16) Acetone	3.78	43	60362	57.700	ug/l	98
18) Methyl Acetate	4.28	43	4892m	1.946	ug/l	
20) Methylene Chloride	4.50	84	16809	5.598	ug/l	97
43) Isopropyl Acetate	8.13	43	39914	6.641	ug/l	99
79) 2-Chlorotoluene	12.65	91	28193	3.268	ug/l	99
82) 4-Chlorotoluene	12.75	91	31312	3.510	ug/l	98

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

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