

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN063020\
 Data File : VN062276.D
 Acq On : 1 Jul 2020 7:08
 Operator : JC/MD
 Sample : L3153-29
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 51 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 SP2-G

Quant Time: Jul 01 07:55:20 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	204071	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	420813	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	414023	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	147277	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	160200	55.94	ug/l	0.00
Spiked Amount 50.000			Recovery	=	111.88%	
35) Dibromofluoromethane	7.55	113	128376	47.55	ug/l	0.00
Spiked Amount 50.000			Recovery	=	95.10%	
50) Toluene-d8	10.06	98	539665	50.01	ug/l	0.00
Spiked Amount 50.000			Recovery	=	100.02%	
62) 4-Bromofluorobenzene	12.38	95	172504	47.31	ug/l	0.00
Spiked Amount 50.000			Recovery	=	94.62%	
Target Compounds						
16) Acetone	3.78	43	67225	71.899	ug/l	94
18) Methyl Acetate	4.28	43	7243	3.182	ug/l	94
20) Methylene Chloride	4.50	84	27616	10.333	ug/l	97
43) Isopropyl Acetate	8.13	43	34327	6.012	ug/l	98

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN063020\
Data File : VN062276.D
Acq On : 1 Jul 2020 7:08
Operator : JC/MD
Sample : L3153-29
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 51 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
SP2-G

Quant Time: Jul 01 07:55:20 2020
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N062420W.M
Quant Title : SW846 8260
QLast Update : Wed Jun 24 15:33:00 2020
Response via : Initial Calibration

