

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN070920\
 Data File : VN062478.D
 Acq On : 10 Jul 2020 00:40
 Operator : JC/MD
 Sample : L3203-03
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 35 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 RO-24

Quant Time: Jul 10 02:40:14 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N070820W.M
 Quant Title : SW846 8260
 QLast Update : Fri Jul 10 00:55:52 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.63	168	193742	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	360585	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	351088	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	130778	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	152646	51.22	ug/l	0.00
Spiked Amount 50.000			Recovery	=	102.44%	
35) Dibromofluoromethane	7.55	113	112178	50.33	ug/l	0.00
Spiked Amount 50.000			Recovery	=	100.66%	
50) Toluene-d8	10.06	98	469381	53.34	ug/l	0.00
Spiked Amount 50.000			Recovery	=	106.68%	
62) 4-Bromofluorobenzene	12.38	95	169570	52.74	ug/l	0.00
Spiked Amount 50.000			Recovery	=	105.48%	
Target Compounds						
16) Acetone	3.78	43	22427	17.726	ug/l	98
18) Methyl Acetate	4.28	43	5868	1.859	ug/l #	91
20) Methylene Chloride	4.50	84	15760	5.826	ug/l	88
43) Isopropyl Acetate	8.13	43	78396	11.683	ug/l #	92

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN070920\
Data File : VN062478.D
Acq On : 10 Jul 2020 00:40
Operator : JC/MD
Sample : L3203-03
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 35 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
RO-24

Quant Time: Jul 10 02:40:14 2020
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N070820W.M
Quant Title : SW846 8260
QLast Update : Fri Jul 10 00:55:52 2020
Response via : Initial Calibration

