

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN031920\
 Data File : VN060641.D
 Acq On : 19 Mar 2020 16:54
 Operator : JC/MD
 Sample : L1978-12
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 TP-7A

Quant Time: Mar 20 08:13:44 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N031820W.M
 Quant Title : SW846 8260
 QLast Update : Wed Mar 18 08:49:09 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.64	168	199164	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.57	114	329295	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.40	117	313690	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.34	152	142270	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.00	65	129678	47.43	ug/l	0.00
Spiked Amount 50.000			Recovery	=	94.86%	
35) Dibromofluoromethane	7.57	113	98540	49.49	ug/l	0.00
Spiked Amount 50.000			Recovery	=	98.98%	
50) Toluene-d8	10.08	98	409744	50.38	ug/l	0.00
Spiked Amount 50.000			Recovery	=	100.76%	
62) 4-Bromofluorobenzene	12.40	95	152402	51.13	ug/l	0.00
Spiked Amount 50.000			Recovery	=	102.26%	
Target Compounds						
16) Acetone	3.79	43	11571	11.948	ug/l	96
20) Methylene Chloride	4.52	84	7533	3.162	ug/l	96
43) Isopropyl Acetate	8.14	43	52197	9.186	ug/l #	91
44) Trichloroethene	8.82	130	11913	4.910	ug/l	98

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN031920\
Data File : VN060641.D
Acq On : 19 Mar 2020 16:54
Operator : JC/MD
Sample : L1978-12
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
TP-7A

Quant Time: Mar 20 08:13:44 2020
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N031820W.M
Quant Title : SW846 8260
QLast Update : Wed Mar 18 08:49:09 2020
Response via : Initial Calibration

