

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN032420\
 Data File : VN060684.D
 Acq On : 24 Mar 2020 19:46
 Operator : JC/MD
 Sample : L2011-03
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 GHD1-WC-S-SB-0203

Quant Time: Mar 25 09:34:09 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	183704	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.57	114	309163	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.40	117	294210	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.34	152	139577	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.00	65	123831	49.10	ug/l	0.00
Spiked Amount 50.000			Recovery	=	98.20%	
35) Dibromofluoromethane	7.56	113	95007	50.82	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.64%	
50) Toluene-d8	10.08	98	386896	50.67	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.34%	
62) 4-Bromofluorobenzene	12.40	95	145348	51.94	ug/l	0.00
Spiked Amount 50.000			Recovery	=	103.88%	
Target Compounds						
16) Acetone	3.79	43	14372	16.089	ug/l	95
18) Methyl Acetate	4.30	43	2662	1.059	ug/l #	68
20) Methylene Chloride	4.52	84	5246	2.387	ug/l	92
25) 2-Butanone	6.80	43	3407	2.357	ug/l	94
43) Isopropyl Acetate	8.14	43	55214	10.350	ug/l #	89

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN032420\
Data File : VN060684.D
Acq On : 24 Mar 2020 19:46
Operator : JC/MD
Sample : L2011-03
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 24 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
GHD1-WC-S-SB-0203

Quant Time: Mar 25 09:34:09 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

