

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN021720\
 Data File : VN060124.D
 Acq On : 17 Feb 2020 17:00
 Operator : JC/MD
 Sample : L1530-02
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 IDW-LGAC-200213

Quant Time: Feb 18 02:42:00 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N021220W.M
 Quant Title : SW846 8260
 QLast Update : Wed Feb 12 14:12:21 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.64	168	181294	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.57	114	288633	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.40	117	267367	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.34	152	121727	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.00	65	111206	50.41	ug/l	0.00
Spiked Amount 50.000			Recovery	=	100.82%	
35) Dibromofluoromethane	7.57	113	87673	49.41	ug/l	0.00
Spiked Amount 50.000			Recovery	=	98.82%	
50) Toluene-d8	10.08	98	350896	50.22	ug/l	0.00
Spiked Amount 50.000			Recovery	=	100.44%	
62) 4-Bromofluorobenzene	12.40	95	126105	49.12	ug/l	0.00
Spiked Amount 50.000			Recovery	=	98.24%	
Target Compounds						
16) Acetone	3.78	43	50256	52.227	ug/l	99
20) Methylene Chloride	4.52	84	3632	1.892	ug/l #	89
25) 2-Butanone	6.81	43	7521	6.215	ug/l	97
43) Isopropyl Acetate	8.14	43	17841	4.320	ug/l #	88

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN021720\
Data File : VN060124.D
Acq On : 17 Feb 2020 17:00
Operator : JC/MD
Sample : L1530-02
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
IDW-LGAC-200213

Quant Time: Feb 18 02:42:00 2020
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N021220W.M
Quant Title : SW846 8260
QLast Update : Wed Feb 12 14:12:21 2020
Response via : Initial Calibration

