

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN042220\
 Data File : VN061136.D
 Acq On : 22 Apr 2020 18:46
 Operator : JC/MD
 Sample : L2330-10
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 FB-SO-2020-04-16

Quant Time: Apr 23 08:10:32 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N042220W.M
 Quant Title : SW846 8260
 QLast Update : Wed Apr 22 15:26:40 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.40	168	196112	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.35	114	325531	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.21	117	305455	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.15	152	149108	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.77	65	120734	46.43	ug/l	0.00
Spiked Amount 50.000			Recovery	=	92.86%	
35) Dibromofluoromethane	7.31	113	67403	36.28	ug/l	0.00
Spiked Amount 50.000			Recovery	=	72.56%	
50) Toluene-d8	9.88	98	396299	49.78	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.56%	
62) 4-Bromofluorobenzene	12.21	95	151130	50.69	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.38%	
Target Compounds						
16) Acetone	3.55	43	29740	20.772	ug/l	98
25) 2-Butanone	6.52	43	8132	5.739	ug/l	91
29) Tetrahydrofuran	6.91	42	2523	2.586	ug/l	96
51) 4-Methyl-2-Pentanone	9.78	43	10874	3.440	ug/l	96

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN042220\
Data File : VN061136.D
Acq On : 22 Apr 2020 18:46
Operator : JC/MD
Sample : L2330-10
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
FB-SO-2020-04-16

Quant Time: Apr 23 08:10:32 2020
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N042220W.M
Quant Title : SW846 8260
QLast Update : Wed Apr 22 15:26:40 2020
Response via : Initial Calibration

