

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN021720\
 Data File : VN060123.D
 Acq On : 17 Feb 2020 16:38
 Operator : JC/MD
 Sample : L1530-01
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 IDW-MCM-200213

Quant Time: Feb 18 02:39:51 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	181148	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.57	114	293347	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.40	117	282451	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.34	152	127841	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.00	65	114639	52.01	ug/l	0.00
Spiked Amount 50.000			Recovery	=	104.02%	
35) Dibromofluoromethane	7.57	113	88342	48.99	ug/l	0.00
Spiked Amount 50.000			Recovery	=	97.98%	
50) Toluene-d8	10.08	98	361850	50.96	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.92%	
62) 4-Bromofluorobenzene	12.40	95	133991	51.35	ug/l	0.00
Spiked Amount 50.000			Recovery	=	102.70%	
Target Compounds						
16) Acetone	3.78	43	31694	32.963	ug/l	100
18) Methyl Acetate	4.30	43	3462	1.917	ug/l #	66
20) Methylene Chloride	4.52	84	9663	5.037	ug/l	94
25) 2-Butanone	6.81	43	8582	7.097	ug/l #	79
43) Isopropyl Acetate	8.14	43	114786	27.345	ug/l #	89
65) Chlorobenzene	11.43	112	256446	44.564	ug/l	100

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

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